

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssptanscl625

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS	1		Web Page for STN Seminar Schedule - N. America
NEWS	2	OCT 02	CA/CAPlus enhanced with pre-1907 records from Chemisches Zentralblatt
NEWS	3	OCT 19	BEILSTEIN updated with new compounds
NEWS	4	NOV 15	Derwent Indian patent publication number format enhanced
NEWS	5	NOV 19	WPIX enhanced with XML display format
NEWS	6	NOV 30	ICSD reloaded with enhancements
NEWS	7	DEC 04	LINPADOCDB now available on STN
NEWS	8	DEC 14	BEILSTEIN pricing structure to change
NEWS	9	DEC 17	USPATOLD added to additional database clusters
NEWS	10	DEC 17	IMSDRUGCONF removed from database clusters and STN
NEWS	11	DEC 17	DGENE now includes more than 10 million sequences
NEWS	12	DEC 17	TOXCENTER enhanced with 2008 MeSH vocabulary in MEDLINE segment
NEWS	13	DEC 17	MEDLINE and LMEDLINE updated with 2008 MeSH vocabulary
NEWS	14	DEC 17	CA/CAPlus enhanced with new custom IPC display formats
NEWS	15	DEC 17	STN Viewer enhanced with full-text patent content from USPATOLD
NEWS	16	JAN 02	STN pricing information for 2008 now available
NEWS	17	JAN 16	CAS patent coverage enhanced to include exemplified prophetic substances
NEWS	18	JAN 28	USPATFULL, USPAT2, and USPATOLD enhanced with new custom IPC display formats
NEWS	19	JAN 28	MARPAT searching enhanced
NEWS	20	JAN 28	USGENE now provides USPTO sequence data within 3 days of publication
NEWS	21	JAN 28	TOXCENTER enhanced with reloaded MEDLINE segment
NEWS	22	JAN 28	MEDLINE and LMEDLINE reloaded with enhancements
NEWS	23	FEB 08	STN Express, Version 8.3, now available
NEWS	24	FEB 20	PCI now available as a replacement to DPCI
NEWS	25	FEB 25	IFIREF reloaded with enhancements
NEWS	26	FEB 25	IMSPRODUCT reloaded with enhancements
NEWS	27	FEB 29	WPINDEX/WPIDS/WPIX enhanced with ECLA and current U.S. National Patent Classification
NEWS	28	MAR 31	IFICDB, IFIPAT, and IFIUIDB enhanced with new custom IPC display formats
NEWS	29	MAR 31	CAS REGISTRY enhanced with additional experimental spectra
NEWS	30	MAR 31	CA/CAPlus and CASREACT patent number format for U.S. applications updated
NEWS	31	MAR 31	LPCI now available as a replacement to LDPCI
NEWS	32	MAR 31	EMBASE, EMBAL, and LEMBASE reloaded with enhancements

NEWS EXPRESS FEBRUARY 08 CURRENT WINDOWS VERSION IS V8.3,
AND CURRENT DISCOVER FILE IS DATED 20 FEBRUARY 2008

NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS LOGIN Welcome Banner and News Items
NEWS IPC8 For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that specific topic.

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* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 16:54:29 ON 03 APR 2008

=> fil reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 16:54:48 ON 03 APR 2008

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STRUCTURE FILE UPDATES: 2 APR 2008 HIGHEST RN 1011757-08-6
DICTIONARY FILE UPDATES: 2 APR 2008 HIGHEST RN 1011757-08-6

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

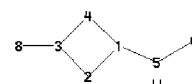
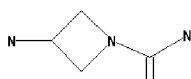
Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10584435.str



```

chain nodes :
5 7
ring nodes :
1 2 3 4
ring/chain nodes :
6 8
chain bonds :
1-5 3-8 5-6 5-7
ring bonds :
1-2 1-4 2-3 3-4
exact/norm bonds :
1-2 1-4 1-5 2-3 3-4 3-8 5-6 5-7

```

G1:O,S,N

Match level :

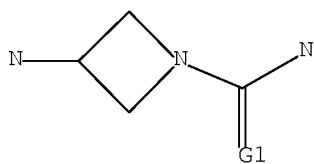
1:Atom 2:Atom 3:Atom 4:Atom 5:CLASS 6:CLASS 7:CLASS 8:CLASS

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



G1 O,S,N

Structure attributes must be viewed using STN Express query preparation.

=> s sss sam l1

SAMPLE SEARCH INITIATED 16:55:07 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 546 TO ITERATE

100.0% PROCESSED 546 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

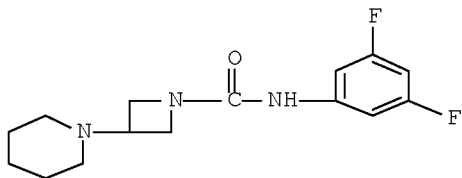
50 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**
PROJECTED ITERATIONS: 9519 TO 12321
PROJECTED ANSWERS: 1830 TO 3170

L2 50 SEA SSS SAM L1

=> d scan

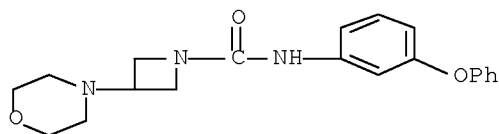
L2 50 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-(1-piperidinyl)-
MF C15 H19 F2 N3 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

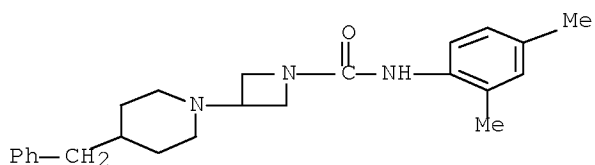
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):3

L2 50 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN 1-Azetidinecarboxamide, 3-(4-morpholinyl)-N-(3-phenoxyphenyl)-
MF C20 H23 N3 O3



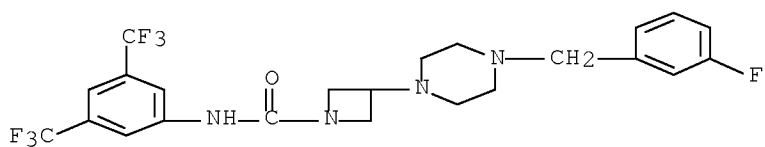
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 50 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]-
MF C24 H31 N3 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 50 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-fluorophenyl)methyl]-1-piperazinyl]-
 MF C23 H23 F7 N4 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> s sss full l1

FULL SEARCH INITIATED 16:55:38 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 10248 TO ITERATE

100.0% PROCESSED 10248 ITERATIONS

2198 ANSWERS

SEARCH TIME: 00.00.01

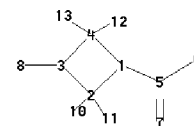
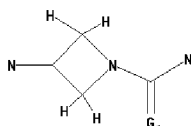
L3 2198 SEA SSS FUL L1

=> save l3 haba10584435/a

ANSWER SET L3 HAS BEEN SAVED AS 'HABA10584435/A'

=>

Uploading C:\Program Files\Stnexp\Queries\10584435A.str



```

chain nodes :
5 7 10 11 12 13
ring nodes :
1 2 3 4
ring/chain nodes :
6 8
chain bonds :
1-5 2-10 2-11 3-8 4-12 4-13 5-6 5-7
ring bonds :
1-2 1-4 2-3 3-4
exact/norm bonds :
1-2 1-4 1-5 2-3 3-4 3-8 5-6 5-7
exact bonds :
2-10 2-11 4-12 4-13

```

G1:O,S,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:CLASS 6:CLASS 7:CLASS 8:CLASS 10:CLASS
11:CLASS 12:CLASS 13:CLASS

L4 STRUCTURE UPLOADED

=> s sss l4 sam subset=l3

SAMPLE SUBSET SEARCH INITIATED 16:58:38 FILE 'REGISTRY'

SAMPLE SUBSET SCREEN SEARCH COMPLETED - 125 TO ITERATE

100.0% PROCESSED 125 ITERATIONS

50 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

PROJECTIONS (WITHIN SPECIFIED SUBSET):

ONLINE **COMPLETE**

PROJECTED ITERATIONS (WITHIN SPECIFIED SUBSET):

1830 TO 3170

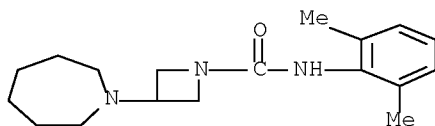
PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET):

915 TO 1925

L5 50 SEA SUB=L3 SSS SAM L4

=> d scan

L5 50 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-(hexahydro-1H-azepin-1-yl)-
MF C18 H27 N3 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> s sss l4 full subset=l3

FULL SUBSET SEARCH INITIATED 16:58:54 FILE 'REGISTRY'

FULL SUBSET SCREEN SEARCH COMPLETED - 2198 TO ITERATE

100.0% PROCESSED 2198 ITERATIONS

1166 ANSWERS

SEARCH TIME: 00.00.01

L6 1166 SEA SUB=L3 SSS FUL L4

=> fil caplu

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

223.22

223.43

FILE 'CAPLUS' ENTERED AT 16:59:02 ON 03 APR 2008

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FILE COVERS 1907 - 3 Apr 2008 VOL 148 ISS 14

FILE LAST UPDATED: 2 Apr 2008 (20080402/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/infopolicy.html>

=> s l6

L7 34 L6

```
=> e us 2006-584435/apps
E1          9      US2006-584429/AP
E2          1      US2006-584432/AP
E3          2 -->  US2006-584435/AP
E4          0      US2006-584435/PRN
E5          1      US2006-584436/AP
E6          1      US2006-584439/AP
E7          1      US2006-584440/AP
E8          1      US2006-584444/AP
E9          1      US2006-584445/AP
E10         2      US2006-584449/AP
E11         2      US2006-584450/AP
E12         1      US2006-584453/AP
```

```
=> s e3
L8          2 US2006-584435/AP
```

```
=> sel rn
E1 THROUGH E927 ASSIGNED
```

```
=> fil reg
COST IN U.S. DOLLARS                SINCE FILE      TOTAL
                                   ENTRY      SESSION
FULL ESTIMATED COST                3.65      227.08
```

FILE 'REGISTRY' ENTERED AT 17:00:33 ON 03 APR 2008
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```
STRUCTURE FILE UPDATES:    2 APR 2008  HIGHEST RN 1011757-08-6
DICTIONARY FILE UPDATES:  2 APR 2008  HIGHEST RN 1011757-08-6
```

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

Please note that search-term pricing does apply when
 conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
 predicted properties as well as tags indicating availability of
 experimental property data in the original document. For information
 on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

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1 182372-13-0/BI
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=> s 19 and 13

L10 917 L9 AND L3

=> d his

(FILE 'HOME' ENTERED AT 16:54:29 ON 03 APR 2008)

FILE 'REGISTRY' ENTERED AT 16:54:48 ON 03 APR 2008

L1 STRUCTURE UPLOADED
L2 50 S SSS SAM L1
L3 2198 S SSS FULL L1
SAVE L3 HABA10584435/A
L4 STRUCTURE UPLOADED
L5 50 S SSS L4 SAM SUB=L3
L6 1166 S SSS L4 FULL SUB=L3

FILE 'CAPLUS' ENTERED AT 16:59:02 ON 03 APR 2008

L7 34 S L6
E US 2006-584435/APPS
L8 2 S E3
SEL RN

FILE 'REGISTRY' ENTERED AT 17:00:33 ON 03 APR 2008

L9 927 S E1-E927
L10 917 S L9 AND L3

=> s 19 and 16

L11 917 L9 AND L6

=> fil caplu

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	2.76	229.84

FILE 'CAPLUS' ENTERED AT 17:04:01 ON 03 APR 2008

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FILE COVERS 1907 - 3 Apr 2008 VOL 148 ISS 14

FILE LAST UPDATED: 2 Apr 2008 (20080402/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/infopolicy.html>

=> s l11

L12 2 L11

=> s l2 and l8

3 L2

L13 1 L2 AND L8

=> fil reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

2.40

232.24

FILE 'REGISTRY' ENTERED AT 17:07:13 ON 03 APR 2008

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STRUCTURE FILE UPDATES: 2 APR 2008 HIGHEST RN 1011757-08-6

DICTIONARY FILE UPDATES: 2 APR 2008 HIGHEST RN 1011757-08-6

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TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stdoc/properties.html>

=> d his

(FILE 'HOME' ENTERED AT 16:54:29 ON 03 APR 2008)

FILE 'REGISTRY' ENTERED AT 16:54:48 ON 03 APR 2008

L1 STRUCTURE UPLOADED

L2 50 S SSS SAM L1

L3 2198 S SSS FULL L1

SAVE L3 HABA10584435/A

L4 STRUCTURE UPLOADED

L5 50 S SSS L4 SAM SUB=L3

L6 1166 S SSS L4 FULL SUB=L3

FILE 'CAPLUS' ENTERED AT 16:59:02 ON 03 APR 2008

L7 34 S L6

E US 2006-584435/APPS

L8 2 S E3

SEL RN

FILE 'REGISTRY' ENTERED AT 17:00:33 ON 03 APR 2008

L9 927 S E1-E927

L10 917 S L9 AND L3

L11 917 S L9 AND L6

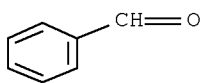
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L12 2 S L11
L13 1 S L2 AND L8

FILE 'REGISTRY' ENTERED AT 17:07:13 ON 03 APR 2008

=> s l9 not l11
L14 10 L9 NOT L11

=> d scan

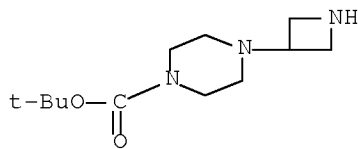
L14 10 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Benzaldehyde
MF C7 H6 O
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):10

L14 10 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN 1-Piperazinecarboxylic acid, 4-(3-azetidiny)-, 1,1-dimethylethyl ester
MF C12 H23 N3 O2
CI COM



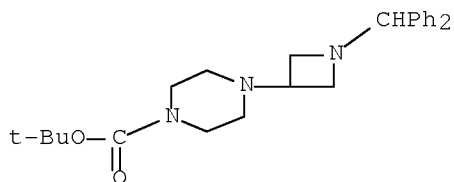
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 10 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Kinase (phosphorylating), RhoA-assocd. coiled-coil forming protein
MF Unspecified
CI MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 10 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN 1-Piperazinecarboxylic acid, 4-[1-(diphenylmethyl)-3-azetidiny]-,
 1,1-dimethylethyl ester
 MF C25 H33 N3 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 10 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Endothelin
 MF Unspecified
 CI MAN

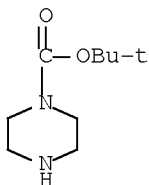
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

L14 10 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Synthetase, thromboxane
 MF Unspecified
 CI MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

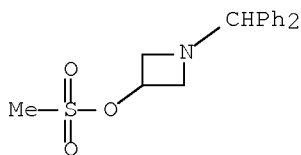
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 10 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN 1-Piperazinecarboxylic acid, 1,1-dimethylethyl ester
 MF C9 H18 N2 O2
 CI COM



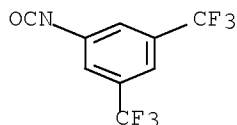
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 10 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN 3-Azetidinol, 1-(diphenylmethyl)-, 3-methanesulfonate
 MF C17 H19 N O3 S
 CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 10 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Benzene, 1-isocyanato-3,5-bis(trifluoromethyl)-
 MF C9 H3 F6 N O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L14 10 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Synthetase, poly(adenosine diphosphoribose)
 MF Unspecified
 CI MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> d his

(FILE 'HOME' ENTERED AT 16:54:29 ON 03 APR 2008)

FILE 'REGISTRY' ENTERED AT 16:54:48 ON 03 APR 2008

L1 STRUCTURE UPLOADED
 L2 50 S SSS SAM L1
 L3 2198 S SSS FULL L1
 SAVE L3 HABA10584435/A
 L4 STRUCTURE UPLOADED
 L5 50 S SSS L4 SAM SUB=L3
 L6 1166 S SSS L4 FULL SUB=L3

FILE 'CAPLUS' ENTERED AT 16:59:02 ON 03 APR 2008

L7 34 S L6
 E US 2006-584435/APPS

L8 2 S E3
 SEL RN

FILE 'REGISTRY' ENTERED AT 17:00:33 ON 03 APR 2008

L9 927 S E1-E927
L10 917 S L9 AND L3
L11 917 S L9 AND L6

FILE 'CAPLUS' ENTERED AT 17:04:01 ON 03 APR 2008

L12 2 S L11
L13 1 S L2 AND L8

FILE 'REGISTRY' ENTERED AT 17:07:13 ON 03 APR 2008

L14 10 S L9 NOT L11

=> s l7 and (ay<2003 or py<2003 or pry<2003)

'2003' NOT A VALID FIELD CODE

'2003' NOT A VALID FIELD CODE

'2003' NOT A VALID FIELD CODE

0 AY<2003

0 PY<2003

0 PRY<2003

L15 0 L7 AND (AY<2003 OR PY<2003 OR PRY<2003)

=> fil caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

1.84

234.08

FILE 'CAPLUS' ENTERED AT 17:09:54 ON 03 APR 2008

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=> d his

(FILE 'HOME' ENTERED AT 16:54:29 ON 03 APR 2008)

FILE 'REGISTRY' ENTERED AT 16:54:48 ON 03 APR 2008

L1 STRUCTURE UPLOADED

L2 50 S SSS SAM L1

L3 2198 S SSS FULL L1
SAVE L3 HABA10584435/A
L4 STRUCTURE UPLOADED
L5 50 S SSS L4 SAM SUB=L3
L6 1166 S SSS L4 FULL SUB=L3

FILE 'CAPLUS' ENTERED AT 16:59:02 ON 03 APR 2008
L7 34 S L6
E US 2006-584435/APPS
L8 2 S E3
SEL RN

FILE 'REGISTRY' ENTERED AT 17:00:33 ON 03 APR 2008
L9 927 S E1-E927
L10 917 S L9 AND L3
L11 917 S L9 AND L6

FILE 'CAPLUS' ENTERED AT 17:04:01 ON 03 APR 2008
L12 2 S L11
L13 1 S L2 AND L8

FILE 'REGISTRY' ENTERED AT 17:07:13 ON 03 APR 2008
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L15 0 S L7 AND (AY<2003 OR PY<2003 OR PRY<2003)

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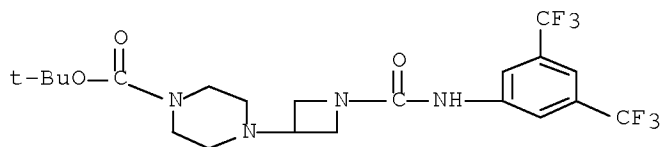
L17 ANSWER 1 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN
ACCESSION NUMBER: 2005:612249 CAPLUS Full-text
DOCUMENT NUMBER: 143:133264
TITLE: Preparation of azetidine ring compounds as EDG-5
antagonists
INVENTOR(S): Habashita, Hiromu; Kurata, Haruto; Nakade, Shinji;
Ono, Takeji
PATENT ASSIGNEE(S): Ono Pharmaceutical Co., Ltd., Japan
SOURCE: PCT Int. Appl., 194 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 2
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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WO 2005063704	A1	20050714	WO 2004-JP19400	20041224 <--
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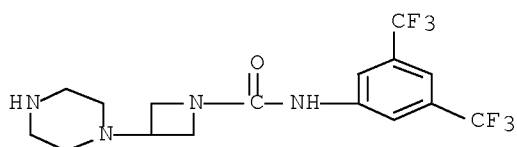
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 LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI,
 NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY,
 TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
 RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM,
 AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK,
 EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC, NL, PL, PT,
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 EP 1698375 A1 20060906 EP 2004-807756 20041224 <--
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 PRIORITY APPLN. INFO.: JP 2003-429948 A 20031225 <--
 JP 2002-185546 A 20020626 <--
 WO 2004-JP19400 W 20041224
 OTHER SOURCE(S): MARPAT 143:133264
 GI

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB Title compds. I [ring A = (un)substituted azetidine ring; X = O, etc.; R1, R2,
 R3, R4 = H, (un)substituted hydrocarbon, etc.] were prepared For example,
 hydrogenolysis of tert-Bu 4-[1-(diphenylmethyl)azetidin-3- yl]piperazine-1-
 carboxylate, e.g., prepared from tert-Bu piperazine-1-carboxylate, using
 Pd(OH)₂ followed by treatment with 1-isocyanato-3,5-
 bis(trifluoromethyl)benzene afforded compound II. In EDG-5 (endothelial
 differentiation gene 5) antagonistic assays, compound III exhibited the IC₅₀
 value of 950 nM. Compds. I are claimed useful for the treatment of
 hypertension, myocardial infarction, etc. Formulations are given.
 IT 858648-34-7P 858656-77-6P
 RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic
 preparation); THU (Therapeutic use); BIOL (Biological study); PREP
 (Preparation); RACT (Reactant or reagent); USES (Uses)
 (preparation of azetidine ring compds. as EDG-5 antagonists for treatment
 of
 hypertension, myocardial infarction, etc.)
 RN 858648-34-7 CAPLUS
 CN 1-Piperazinecarboxylic acid, 4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]
 carbonyl]-3-azetidinyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)



RN 858656-77-6 CAPLUS
 CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(1-
 piperazinyl)- (CA INDEX NAME)



IT 642495-73-6P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide
 642498-98-4P, N-(3,5-Dimethylphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide
 642498-99-5P, 3-[4-[4-(Trifluoromethoxy)phenyl]piperazin-1-yl]-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858648-35-8P
 858648-36-9P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(2-methoxyphenyl)piperazin-1-yl]azetidine-1-carboxamide 858648-37-0P
 , N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(3-methoxyphenyl)piperazin-1-yl]azetidine-1-carboxamide 858648-38-1P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(4-methoxyphenyl)piperazin-1-yl]azetidine-1-carboxamide 858648-39-2P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-[3-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-40-5P, N-(2-Ethylphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858648-41-6P, N-(2,4-Dimethylphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858648-42-7P, 3-(4-Methylpiperazin-1-yl)-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858648-43-8P, N-(3,5-Dimethylphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858648-44-9P, Methyl 3-([3-(4-methylpiperazin-1-yl)azetidin-1-yl]carbonylamino)benzoate 858648-45-0P, N-(2,4-Dimethylphenyl)-3-pyrrolidin-1-ylazetidine-1-carboxamide 858648-46-1P, N-(3,5-Dichlorophenyl)-3-(4-phenylpiperidin-1-yl)azetidine-1-carboxamide 858648-47-2P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858648-48-3P, N-(3,5-Dichlorophenyl)-3-(2,3-dihydro-1H-indol-1-yl)azetidine-1-carboxamide 858648-49-4P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-(2,3-dihydro-1H-indol-1-yl)azetidine-1-carboxamide 858648-50-7P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858648-51-8P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858648-52-9P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[ethyl(phenyl)amino]azetidine-1-carboxamide 858648-53-0P, N-Phenyl-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-54-1P, N-Butyl-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-55-2P, N-(4-Chlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-56-3P, N-(3-Chlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-57-4P, N-Cyclohexyl-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-58-5P, N-(2-Chlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-59-6P, N-(3,4-Dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-60-9P, N-(2-Methoxyphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-61-0P, N-Hexyl-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-62-1P, N-(3-Methoxyphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-63-2P, N-(4-Methoxyphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-64-3P, 3-[4-[4-(Trifluoromethoxy)phenyl]piperazin-1-yl]-N-

[2-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858648-65-4P,
N-(2,4-Dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-66-5P 858648-67-6P,
N-(2-Fluorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-68-7P, N-(3-Fluorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-69-8P, N-Benzyl-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-70-1P, N-(4-Fluorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-71-2P, 3-[4-[4-(Trifluoromethoxy)phenyl]piperazin-1-yl]-N-[4-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858648-72-3P,
N-(3,5-Dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-73-4P, N-(2,5-Dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-74-5P, N-Pentyl-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-75-6P, N-(2,6-Dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-76-7P, N-(2-Phenylethyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-77-8P, N-(2,3-Dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-78-9P, N-(3-Cyanophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-79-0P, Ethyl 4-[[3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidin-1-yl]carbonyl]amino]benzoate 858648-80-3P, N-(4-Phenoxyphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-81-4P, Ethyl 3-[[3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidin-1-yl]carbonyl]amino]benzoate 858648-82-5P, N-Isopropyl-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-83-6P, N-(3-Phenoxyphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-84-7P, N-(4-Cyanophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-85-8P, N-(3,5-Difluorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-86-9P, N-[3-Fluoro-5-(trifluoromethyl)phenyl]-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-87-0P, N-(3-Chloro-5-fluorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-88-1P, N-[3-(Cyclopentyloxy)phenyl]-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-89-2P, N-[3-(Cyclohexyloxy)phenyl]-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-92-7P, N-(2,6-Dichloropyridin-4-yl)-3-[4-[4-(trifluoromethoxy)phenyl]piperazin-1-yl]azetidine-1-carboxamide 858648-93-8P, 3-(Dimethylamino)-N-hexylazetidine-1-carboxamide 858648-94-9P, 3-(Dimethylamino)-N-(4-fluorophenyl)azetidine-1-carboxamide 858648-95-0P, 3-(Dimethylamino)-N-(3-methylphenyl)azetidine-1-carboxamide 858648-96-1P, 3-(Dimethylamino)-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858648-97-2P, 3-(Dimethylamino)-N-(4-isopropylphenyl)azetidine-1-carboxamide 858648-98-3P, N-(3-Chlorophenyl)-3-(dimethylamino)azetidine-1-carboxamide 858648-99-4P, 3-(Dimethylamino)-N-(2,5-dimethylphenyl)azetidine-1-carboxamide 858649-00-0P, N-(4-Chlorophenyl)-3-(dimethylamino)azetidine-1-carboxamide 858649-01-1P, N-Benzyl-3-(dimethylamino)azetidine-1-carboxamide 858649-02-2P, 3-(Dimethylamino)-N-(1-naphthyl)azetidine-1-carboxamide 858649-03-3P, 3-(Dimethylamino)-N-(2-naphthyl)azetidine-1-carboxamide

858649-04-4P, 3-(Dimethylamino)-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858649-05-5P, 3-(Dimethylamino)-N-(3,4-dimethylphenyl)azetidine-1-carboxamide 858649-06-6P, 3-(Dimethylamino)-N-(4-methylphenyl)azetidine-1-carboxamide 858649-07-7P, N-Cyclohexyl-3-(dimethylamino)azetidine-1-carboxamide 858649-08-8P, 3-(Dimethylamino)-N-(2,6-dimethylphenyl)azetidine-1-carboxamide 858649-09-9P, 3-(Dimethylamino)-N-(2-ethoxyphenyl)azetidine-1-carboxamide 858649-10-2P, 3-(Dimethylamino)-N-(4-ethoxyphenyl)azetidine-1-carboxamide 858649-11-3P, 3-(Dimethylamino)-N-(2-phenylethyl)azetidine-1-carboxamide 858649-12-4P, 3-(Dimethylamino)-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858649-13-5P, N-Cyclopentyl-3-(dimethylamino)azetidine-1-carboxamide 858649-14-6P, 3-(Dimethylamino)-N-(2,4-dimethylphenyl)azetidine-1-carboxamide 858649-15-7P, N-(3,5-Dichlorophenyl)-3-(dimethylamino)azetidine-1-carboxamide 858649-16-8P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-(dimethylamino)azetidine-1-carboxamide 858649-17-9P, 3-(Dimethylamino)-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858649-18-0P, N-(3,5-Difluorophenyl)-3-(dimethylamino)azetidine-1-carboxamide 858649-19-1P, 3-(Dimethylamino)-N-(4-methoxyphenyl)azetidine-1-carboxamide 858649-20-4P, 3-(Dimethylamino)-N-(3,5-dimethylphenyl)azetidine-1-carboxamide 858649-21-5P, 3-(Dimethylamino)-N-(3-fluorophenyl)azetidine-1-carboxamide 858649-22-6P, Methyl 3-([3-(dimethylamino)azetidin-1-yl]carbonyl)amino)benzoate 858649-23-7P, 3-(Dimethylamino)-N-[3-(methylsulfonyl)phenyl]azetidine-1-carboxamide 858649-24-8P, 3-(Diethylamino)-N-propylazetidine-1-carboxamide 858649-25-9P 858649-26-0P, 3-(Diethylamino)-N-hexylazetidine-1-carboxamide 858649-27-1P, 3-(Diethylamino)-N-(4-fluorophenyl)azetidine-1-carboxamide 858649-28-2P, 3-(Diethylamino)-N-(3-methylphenyl)azetidine-1-carboxamide 858649-29-3P, 3-(Diethylamino)-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858649-30-6P, 3-(Diethylamino)-N-(4-isopropylphenyl)azetidine-1-carboxamide 858649-31-7P, N-(3-Chlorophenyl)-3-(diethylamino)azetidine-1-carboxamide 858649-32-8P, 3-(Diethylamino)-N-(2,5-dimethylphenyl)azetidine-1-carboxamide 858649-33-9P, N-(4-Chlorophenyl)-3-(diethylamino)azetidine-1-carboxamide 858649-34-0P, N-Benzyl-3-(diethylamino)azetidine-1-carboxamide 858649-35-1P, 3-(Diethylamino)-N-(1-naphthyl)azetidine-1-carboxamide 858649-36-2P, 3-(Diethylamino)-N-(2-naphthyl)azetidine-1-carboxamide 858649-37-3P, 3-(Diethylamino)-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858649-38-4P, 3-(Diethylamino)-N-(3,4-dimethylphenyl)azetidine-1-carboxamide 858649-39-5P, 3-(Diethylamino)-N-(4-methylphenyl)azetidine-1-carboxamide 858649-40-8P, N-Cyclohexyl-3-(diethylamino)azetidine-1-carboxamide 858649-41-9P, 3-(Diethylamino)-N-(2,6-dimethylphenyl)azetidine-1-carboxamide 858649-42-0P, 3-(Diethylamino)-N-(2-ethoxyphenyl)azetidine-1-carboxamide 858649-43-1P, 3-(Diethylamino)-N-(2-ethylphenyl)azetidine-1-carboxamide 858649-44-2P, 3-(Diethylamino)-N-(4-ethoxyphenyl)azetidine-1-carboxamide 858649-45-3P, 3-(Diethylamino)-N-(2-phenylethyl)azetidine-1-carboxamide 858649-46-4P, 3-(Diethylamino)-N-phenylazetidine-1-carboxamide 858649-47-5P, N-(2-Chlorophenyl)-3-(diethylamino)azetidine-1-carboxamide 858649-48-6P, 3-(Diethylamino)-N-(2-fluorophenyl)azetidine-1-carboxamide 858649-49-7P, 3-(Diethylamino)-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858649-50-0P, N-Cyclopentyl-3-(diethylamino)azetidine-1-carboxamide 858649-51-1P

, 3-(Diethylamino)-N-(2,4-dimethylphenyl)azetidine-1-carboxamide 858649-52-2P, N-(3,5-Dichlorophenyl)-3-(diethylamino)azetidine-1-carboxamide 858649-53-3P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-(diethylamino)azetidine-1-carboxamide 858649-54-4P, 3-(Diethylamino)-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858649-55-5P, 3-(Diethylamino)-N-(3,5-difluorophenyl)azetidine-1-carboxamide 858649-56-6P, 3-(Diethylamino)-N-(4-methoxyphenyl)azetidine-1-carboxamide 858649-57-7P, 3-(Diethylamino)-N-(3,5-dimethylphenyl)azetidine-1-carboxamide 858649-58-8P, 3-(Diethylamino)-N-(3-fluorophenyl)azetidine-1-carboxamide 858649-59-9P, Methyl 3-([(3-(diethylamino)azetidin-1-yl]carbonyl]amino)benzoate 858649-60-2P, 3-(Diethylamino)-N-[3-(methylsulfanyl)phenyl]azetidine-1-carboxamide 858649-61-3P, 3-(Diethylamino)-N-(2-methylphenyl)azetidine-1-carboxamide 858649-62-4P, 3-(Diisopropylamino)-N-propylazetidine-1-carboxamide 858649-63-5P 858649-64-6P, 3-(Diisopropylamino)-N-hexylazetidine-1-carboxamide 858649-65-7P, 3-(Diisopropylamino)-N-(4-fluorophenyl)azetidine-1-carboxamide 858649-66-8P, 3-(Diisopropylamino)-N-(3-methylphenyl)azetidine-1-carboxamide 858649-67-9P, 3-(Diisopropylamino)-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858649-68-0P, 3-(Diisopropylamino)-N-(4-isopropylphenyl)azetidine-1-carboxamide 858649-69-1P, N-(3-Chlorophenyl)-3-(diisopropylamino)azetidine-1-carboxamide 858649-70-4P, 3-(Diisopropylamino)-N-(2,5-dimethylphenyl)azetidine-1-carboxamide 858649-71-5P, N-(4-Chlorophenyl)-3-(diisopropylamino)azetidine-1-carboxamide 858649-72-6P, N-Benzyl-3-(diisopropylamino)azetidine-1-carboxamide 858649-73-7P, 3-(Diisopropylamino)-N-(1-naphthyl)azetidine-1-carboxamide 858649-74-8P, 3-(Diisopropylamino)-N-(2-naphthyl)azetidine-1-carboxamide 858649-75-9P, 3-(Diisopropylamino)-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858649-76-0P, 3-(Diisopropylamino)-N-(3,4-dimethylphenyl)azetidine-1-carboxamide 858649-77-1P, 3-(Diisopropylamino)-N-(4-methylphenyl)azetidine-1-carboxamide 858649-78-2P, N-Cyclohexyl-3-(diisopropylamino)azetidine-1-carboxamide 858649-79-3P, 3-(Diisopropylamino)-N-(2,6-dimethylphenyl)azetidine-1-carboxamide 858649-80-6P, 3-(Diisopropylamino)-N-(2-ethoxyphenyl)azetidine-1-carboxamide 858649-81-7P, 3-(Diisopropylamino)-N-(2-ethylphenyl)azetidine-1-carboxamide 858649-82-8P, 3-(Diisopropylamino)-N-(4-ethoxyphenyl)azetidine-1-carboxamide 858649-83-9P, 3-(Diisopropylamino)-N-(2-phenylethyl)azetidine-1-carboxamide 858649-84-0P, 3-(Diisopropylamino)-N-phenylazetidine-1-carboxamide 858649-85-1P, N-(2-Chlorophenyl)-3-(diisopropylamino)azetidine-1-carboxamide 858649-86-2P, 3-(Diisopropylamino)-N-(2-fluorophenyl)azetidine-1-carboxamide 858649-87-3P, 3-(Diisopropylamino)-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858649-88-4P, N-Cyclopentyl-3-(diisopropylamino)azetidine-1-carboxamide 858649-89-5P, 3-(Diisopropylamino)-N-(2,4-dimethylphenyl)azetidine-1-carboxamide 858649-90-8P, N-(3,5-Dichlorophenyl)-3-(diisopropylamino)azetidine-1-carboxamide 858649-91-9P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-(diisopropylamino)azetidine-1-carboxamide 858649-92-0P, 3-(Diisopropylamino)-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858649-93-1P, N-(3,5-Difluorophenyl)-3-(diisopropylamino)azetidine-1-carboxamide 858649-94-2P, 3-(Diisopropylamino)-N-(4-methoxyphenyl)azetidine-1-carboxamide 858649-95-3P, 3-(Diisopropylamino)-N-(3,5-dimethylphenyl)azetidine-1-carboxamide 858649-96-4P, 3-(Diisopropylamino)-N-(3-fluorophenyl)azetidine-1-carboxamide 858649-97-5P, Methyl 3-([(3-(diisopropylamino)azetidin-1-

yl]carbonyl]amino)benzoate 858649-98-6P, 3-(Diisopropylamino)-N-[3-(methylsulfanyl)phenyl]azetidine-1-carboxamide 858649-99-7P, 3-(Diisopropylamino)-N-(2-methylphenyl)azetidine-1-carboxamide 858650-00-7P, 3-(Dipropylamino)-N-propylazetidine-1-carboxamide 858650-01-8P 858650-02-9P, 3-(Dipropylamino)-N-hexylazetidine-1-carboxamide 858650-03-0P, 3-(Dipropylamino)-N-(4-fluorophenyl)azetidine-1-carboxamide 858650-04-1P, 3-(Dipropylamino)-N-(3-methylphenyl)azetidine-1-carboxamide 858650-05-2P, 3-(Dipropylamino)-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858650-06-3P, 3-(Dipropylamino)-N-(4-isopropylphenyl)azetidine-1-carboxamide 858650-07-4P, N-(3-Chlorophenyl)-3-(dipropylamino)azetidine-1-carboxamide 858650-08-5P, N-(2,5-Dimethylphenyl)-3-(dipropylamino)azetidine-1-carboxamide 858650-09-6P, N-(4-Chlorophenyl)-3-(dipropylamino)azetidine-1-carboxamide 858650-10-9P, N-Benzyl-3-(dipropylamino)azetidine-1-carboxamide 858650-11-0P, 3-(Dipropylamino)-N-(1-naphthyl)azetidine-1-carboxamide 858650-12-1P, 3-(Dipropylamino)-N-(2-naphthyl)azetidine-1-carboxamide 858650-13-2P, 3-(Dipropylamino)-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858650-14-3P, N-(3,4-Dimethylphenyl)-3-(dipropylamino)azetidine-1-carboxamide 858650-15-4P, 3-(Dipropylamino)-N-(4-methylphenyl)azetidine-1-carboxamide 858650-16-5P, N-Cyclohexyl-3-(dipropylamino)azetidine-1-carboxamide 858650-17-6P, N-(2,6-Dimethylphenyl)-3-(dipropylamino)azetidine-1-carboxamide 858650-18-7P, 3-(Dipropylamino)-N-(2-ethoxyphenyl)azetidine-1-carboxamide 858650-19-8P, 3-(Dipropylamino)-N-(2-ethylphenyl)azetidine-1-carboxamide 858650-20-1P, 3-(Dipropylamino)-N-(4-ethoxyphenyl)azetidine-1-carboxamide 858650-21-2P, 3-(Dipropylamino)-N-(2-phenylethyl)azetidine-1-carboxamide 858650-22-3P, 3-(Dipropylamino)-N-phenylazetidine-1-carboxamide 858650-23-4P, N-(2-Chlorophenyl)-3-(dipropylamino)azetidine-1-carboxamide 858650-24-5P, 3-(Dipropylamino)-N-(2-fluorophenyl)azetidine-1-carboxamide 858650-25-6P, 3-(Dipropylamino)-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858650-26-7P, N-Cyclopentyl-3-(dipropylamino)azetidine-1-carboxamide 858650-27-8P, N-(2,4-Dimethylphenyl)-3-(dipropylamino)azetidine-1-carboxamide 858650-28-9P, N-(3,5-Dichlorophenyl)-3-(dipropylamino)azetidine-1-carboxamide 858650-29-0P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-(dipropylamino)azetidine-1-carboxamide 858650-30-3P, 3-(Dipropylamino)-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858650-32-5P, N-(3,5-Difluorophenyl)-3-(Dipropylamino)azetidine-1-carboxamide 858650-34-7P, 3-(Dipropylamino)-N-(4-methoxyphenyl)azetidine-1-carboxamide 858650-36-9P, N-(3,5-Dimethylphenyl)-3-(dipropylamino)azetidine-1-carboxamide 858650-38-1P, 3-(Dipropylamino)-N-(3-fluorophenyl)azetidine-1-carboxamide 858650-40-5P, Methyl 3-([(3-(dipropylamino)azetidin-1-yl]carbonyl]amino)benzoate 858650-42-7P, 3-(Dipropylamino)-N-[3-(methylsulfanyl)phenyl]azetidine-1-carboxamide 858650-44-9P, 3-(Dipropylamino)-N-(2-methylphenyl)azetidine-1-carboxamide 858650-46-1P, 3-[Bis(2-hydroxyethyl)amino]-N-hexylazetidine-1-carboxamide 858650-48-3P, 3-[Bis(2-hydroxyethyl)amino]-N-(4-fluorophenyl)azetidine-1-carboxamide 858650-50-7P, 3-[Bis(2-hydroxyethyl)amino]-N-(3-methylphenyl)azetidine-1-carboxamide 858650-51-8P, 3-[Bis(2-hydroxyethyl)amino]-N-(4-isopropylphenyl)azetidine-1-carboxamide 858650-52-9P, 3-[Bis(2-hydroxyethyl)amino]-N-(3-chlorophenyl)azetidine-1-carboxamide 858650-53-0P, 3-[Bis(2-hydroxyethyl)amino]-N-(2,5-dimethylphenyl)azetidine-1-carboxamide 858650-54-1P,

3-[Bis(2-hydroxyethyl)amino]-N-(4-chlorophenyl)azetidine-1-carboxamide 858650-55-2P, 3-[Bis(2-hydroxyethyl)amino]-N-(1-naphthyl)azetidine-1-carboxamide 858650-56-3P, 3-[Bis(2-hydroxyethyl)amino]-N-(2-naphthyl)azetidine-1-carboxamide 858650-57-4P, 3-[Bis(2-hydroxyethyl)amino]-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858650-58-5P

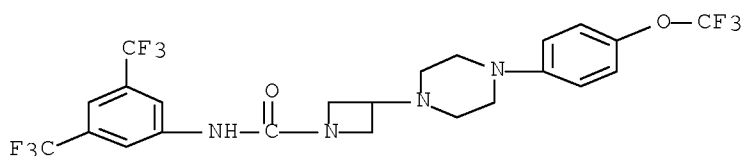
, 3-[Bis(2-hydroxyethyl)amino]-N-(3,4-dimethylphenyl)azetidine-1-carboxamide 858650-59-6P, 3-[Bis(2-hydroxyethyl)amino]-N-(4-methylphenyl)azetidine-1-carboxamide 858650-60-9P, 3-[Bis(2-hydroxyethyl)amino]-N-cyclohexylazetidine-1-carboxamide 858650-61-0P, 3-[Bis(2-hydroxyethyl)amino]-N-(2,6-dimethylphenyl)azetidine-1-carboxamide 858650-62-1P, 3-[Bis(2-hydroxyethyl)amino]-N-(2-ethoxyphenyl)azetidine-1-carboxamide 858650-63-2P, 3-[Bis(2-hydroxyethyl)amino]-N-(2-ethylphenyl)azetidine-1-carboxamide 858650-64-3P, 3-[Bis(2-hydroxyethyl)amino]-N-(4-ethoxyphenyl)azetidine-1-carboxamide 858650-65-4P, 3-[Bis(2-hydroxyethyl)amino]-N-(2-phenylethyl)azetidine-1-carboxamide 858650-66-5P, 3-[Bis(2-hydroxyethyl)amino]-N-phenylazetidine-1-carboxamide 858650-67-6P, 3-[Bis(2-hydroxyethyl)amino]-N-(2-chlorophenyl)azetidine-1-carboxamide 858650-68-7P, 3-[Bis(2-hydroxyethyl)amino]-N-(2-fluorophenyl)azetidine-1-carboxamide 858650-69-8P, 3-[Bis(2-hydroxyethyl)amino]-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858650-70-1P, 3-[Bis(2-hydroxyethyl)amino]-N-cyclopentylazetidine-1-carboxamide 858650-71-2P, 3-[Bis(2-hydroxyethyl)amino]-N-(2,4-dimethylphenyl)azetidine-1-carboxamide 858650-72-3P, 3-[Bis(2-hydroxyethyl)amino]-N-(3,5-dichlorophenyl)azetidine-1-carboxamide 858650-73-4P, 3-[Bis(2-hydroxyethyl)amino]-N-[3,5-bis(trifluoromethyl)phenyl]azetidine-1-carboxamide 858650-74-5P, 3-[Bis(2-hydroxyethyl)amino]-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858650-75-6P, 3-[Bis(2-hydroxyethyl)amino]-N-(3,5-difluorophenyl)azetidine-1-carboxamide 858650-76-7P, 3-[Bis(2-hydroxyethyl)amino]-N-(4-methoxyphenyl)azetidine-1-carboxamide 858650-77-8P, 3-[Bis(2-hydroxyethyl)amino]-N-(3,5-dimethylphenyl)azetidine-1-carboxamide

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of azetidine ring compds. as EDG-5 antagonists for treatment of hypertension, myocardial infarction, etc.)

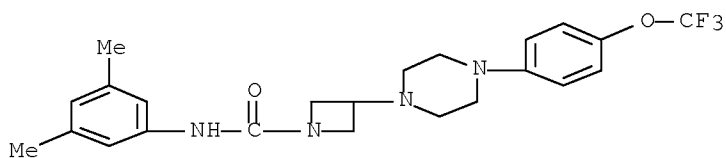
RN 642495-73-6 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



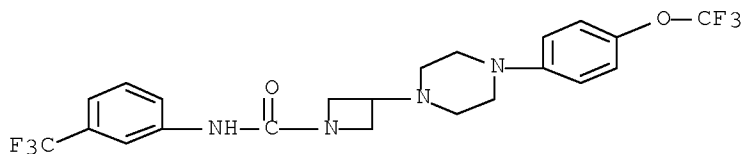
RN 642498-98-4 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



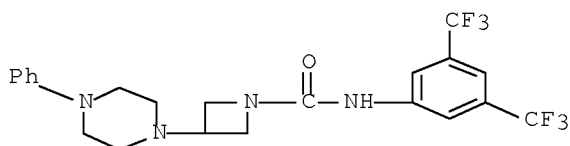
RN 642498-99-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



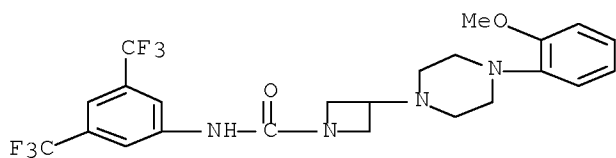
RN 858648-35-8 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)



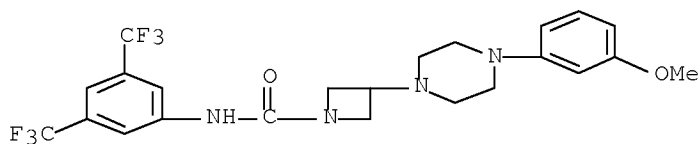
RN 858648-36-9 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(2-methoxyphenyl)-1-piperazinyl]- (CA INDEX NAME)

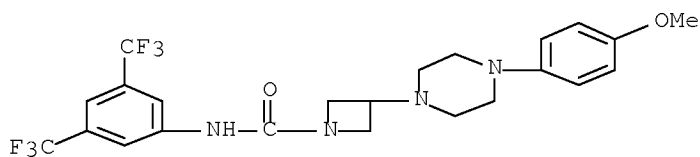


RN 858648-37-0 CAPLUS

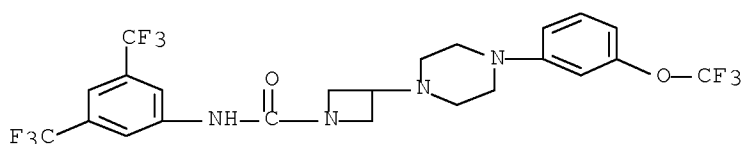
CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(3-methoxyphenyl)-1-piperazinyl]- (CA INDEX NAME)



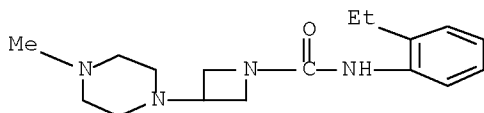
RN 858648-38-1 CAPLUS
 CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(4-methoxyphenyl)-1-piperazinyl]- (CA INDEX NAME)



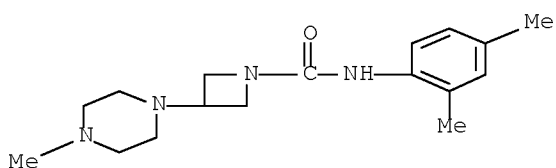
RN 858648-39-2 CAPLUS
 CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[3-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



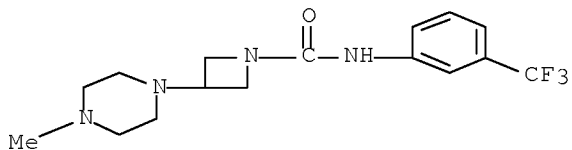
RN 858648-40-5 CAPLUS
 CN 1-Azetidinecarboxamide, N-(2-ethylphenyl)-3-(4-methyl-1-piperazinyl)- (CA INDEX NAME)



RN 858648-41-6 CAPLUS
 CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(4-methyl-1-piperazinyl)- (CA INDEX NAME)

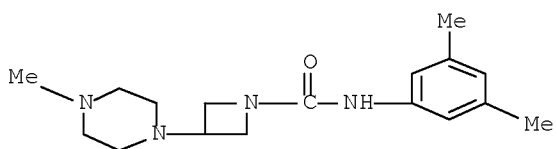


RN 858648-42-7 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(4-methyl-1-piperazinyl)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



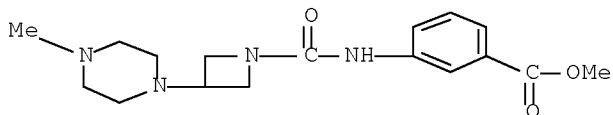
RN 858648-43-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-(4-methyl-1-piperazinyl)-
(CA INDEX NAME)



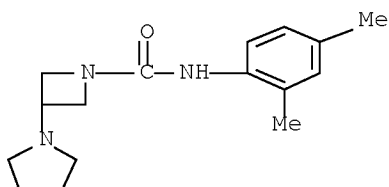
RN 858648-44-9 CAPLUS

CN Benzoic acid, 3-[[[3-(4-methyl-1-piperazinyl)-1-azetidiny]carbonyl]amino]-
, methyl ester (CA INDEX NAME)



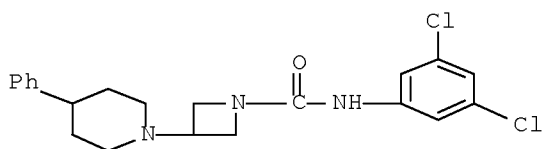
RN 858648-45-0 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(1-pyrrolidinyl)- (CA
INDEX NAME)



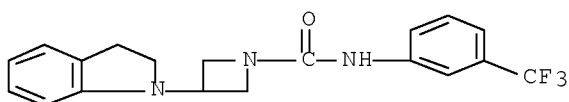
RN 858648-46-1 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(4-phenyl-1-piperidiny)-
(CA INDEX NAME)



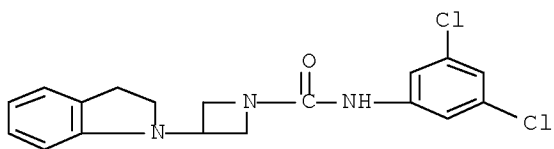
RN 858648-47-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



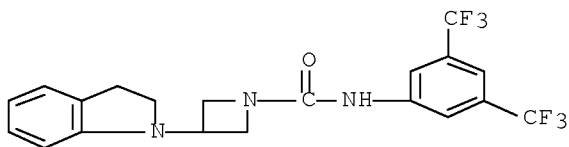
RN 858648-48-3 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(2,3-dihydro-1H-indol-1-yl)- (CA INDEX NAME)



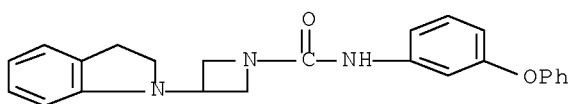
RN 858648-49-4 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(2,3-dihydro-1H-indol-1-yl)- (CA INDEX NAME)

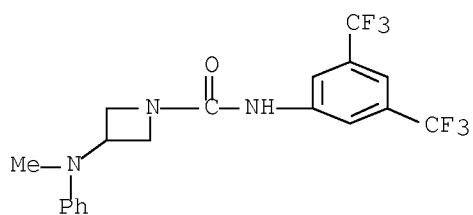


RN 858648-50-7 CAPLUS

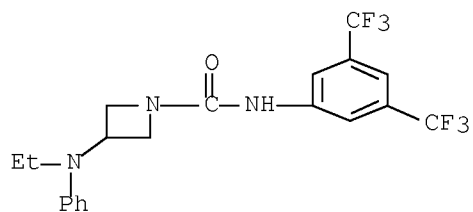
CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(3-phenoxyphenyl)- (CA INDEX NAME)



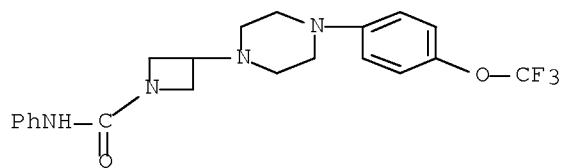
RN 858648-51-8 CAPLUS
 CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(methylphenylamino)- (CA INDEX NAME)



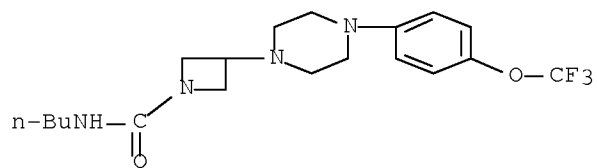
RN 858648-52-9 CAPLUS
 CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(ethylphenylamino)- (CA INDEX NAME)



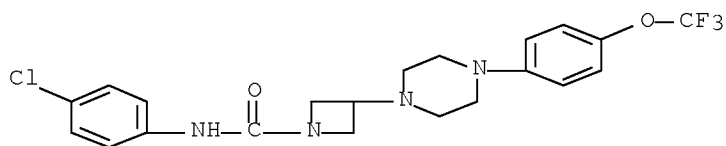
RN 858648-53-0 CAPLUS
 CN 1-Azetidinecarboxamide, N-phenyl-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



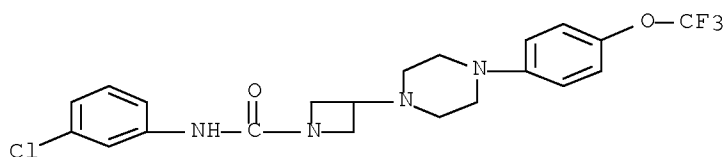
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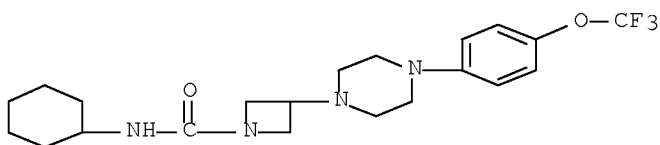
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CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



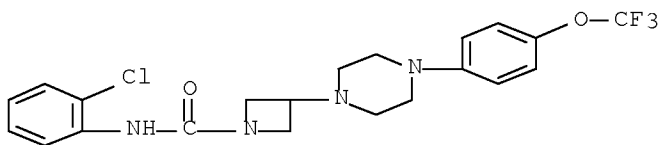
RN 858648-56-3 CAPLUS
CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



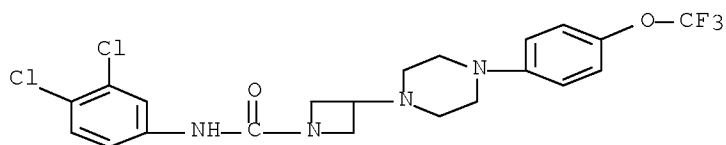
RN 858648-57-4 CAPLUS
CN 1-Azetidinecarboxamide, N-cyclohexyl-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



RN 858648-58-5 CAPLUS
CN 1-Azetidinecarboxamide, N-(2-chlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)

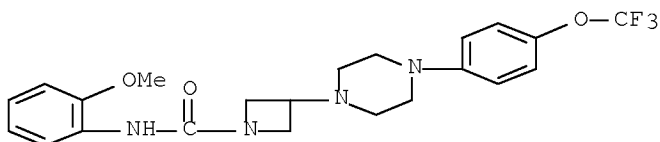


RN 858648-59-6 CAPLUS
CN 1-Azetidinecarboxamide, N-(3,4-dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



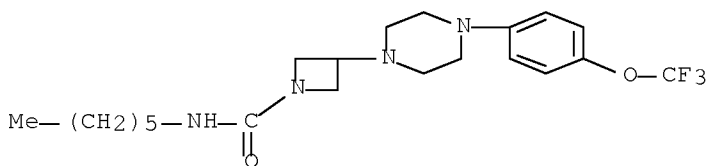
RN 858648-60-9 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-methoxyphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



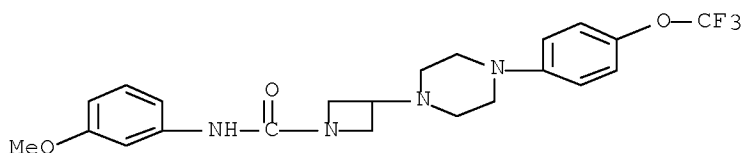
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CN 1-Azetidinecarboxamide, N-hexyl-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



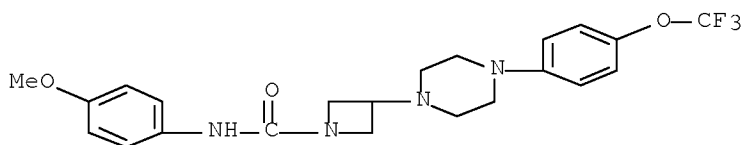
RN 858648-62-1 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-methoxyphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



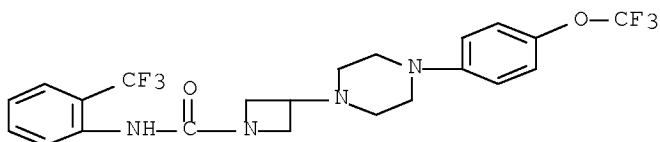
RN 858648-63-2 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-methoxyphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



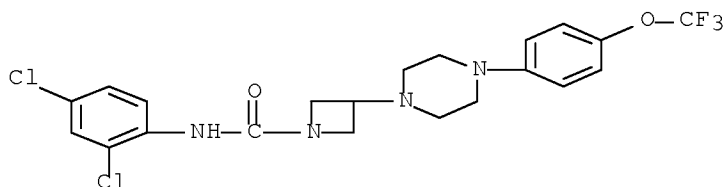
RN 858648-64-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]-N-[2-(trifluoromethyl)phenyl]- (CA INDEX NAME)



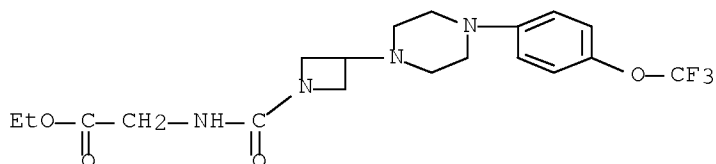
RN 858648-65-4 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,4-dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



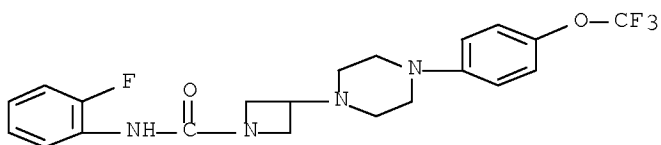
RN 858648-66-5 CAPLUS

CN Glycine, N-[[3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)



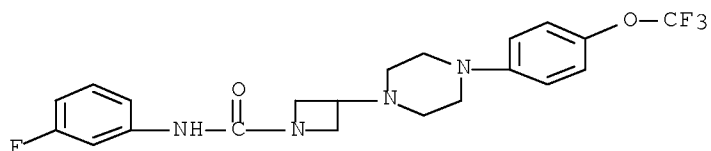
RN 858648-67-6 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-fluorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



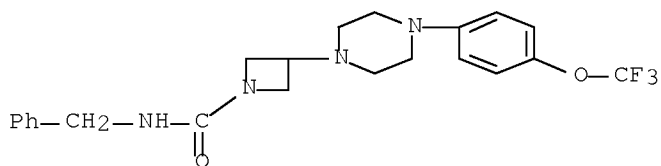
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CN 1-Azetidinecarboxamide, N-(3-fluorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



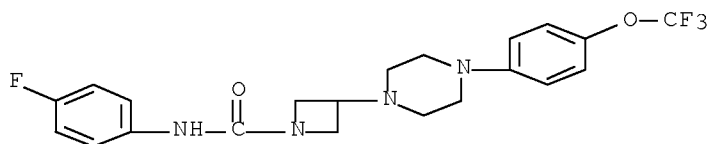
RN 858648-69-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(phenylmethyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



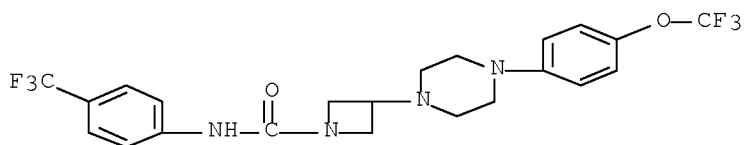
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CN 1-Azetidinecarboxamide, N-(4-fluorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



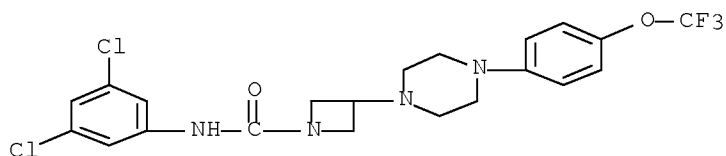
RN 858648-71-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]-N-[4-(trifluoromethyl)phenyl]- (CA INDEX NAME)



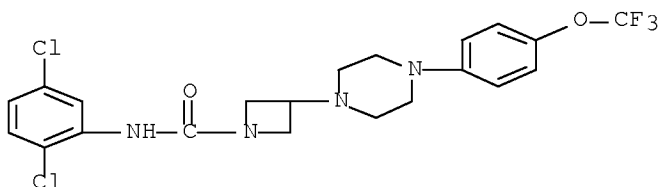
RN 858648-72-3 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



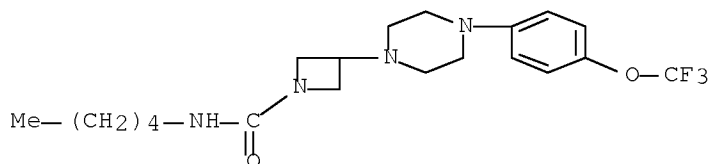
RN 858648-73-4 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,5-dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



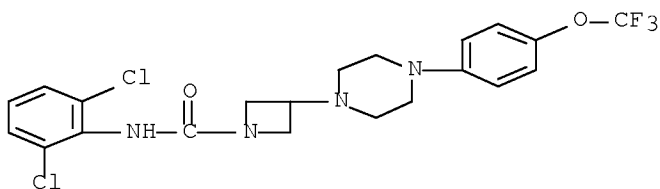
RN 858648-74-5 CAPLUS

CN 1-Azetidinecarboxamide, N-pentyl-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



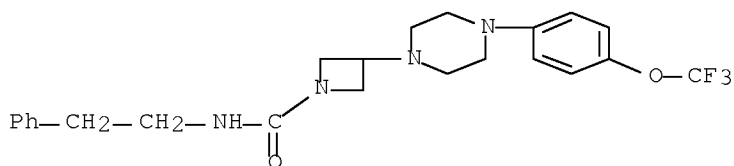
RN 858648-75-6 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,6-dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



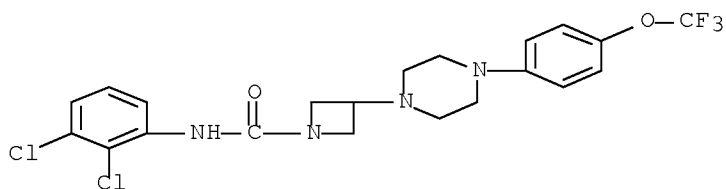
RN 858648-76-7 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-phenylethyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



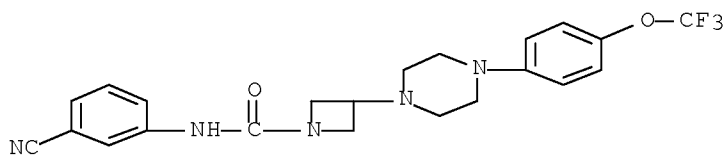
RN 858648-77-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,3-dichlorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



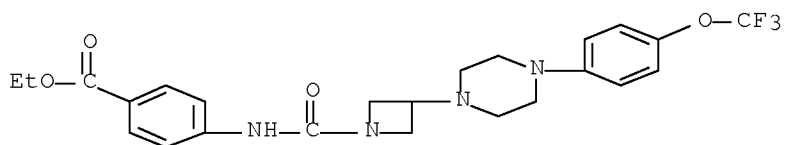
RN 858648-78-9 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-cyanophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



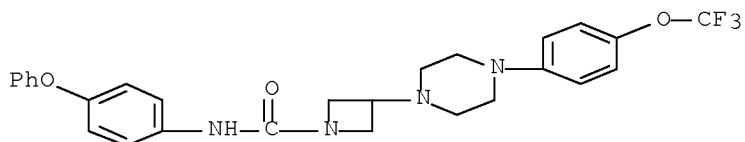
RN 858648-79-0 CAPLUS

CN Benzoic acid, 4-[[[3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]-1-azetidiny]carbonyl]amino]-, ethyl ester (CA INDEX NAME)



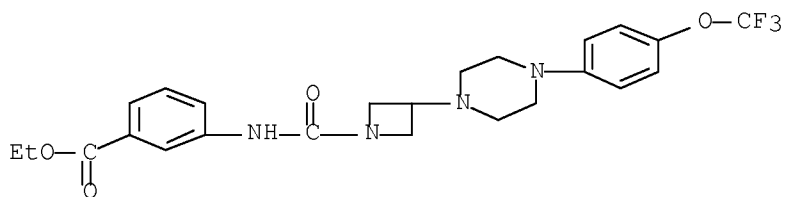
RN 858648-80-3 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-phenoxyphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



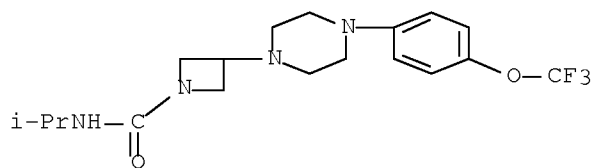
RN 858648-81-4 CAPLUS

CN Benzoic acid, 3-[[[3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]-1-azetidiny]carbonyl]amino]-, ethyl ester (CA INDEX NAME)



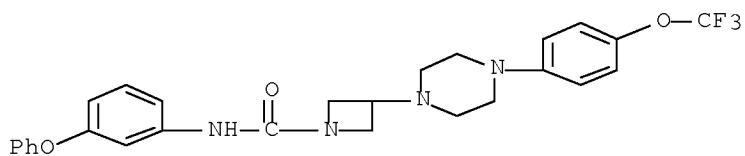
RN 858648-82-5 CAPLUS

CN 1-Azetidinecarboxamide, N-(1-methylethyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



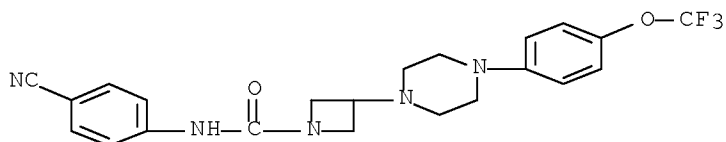
RN 858648-83-6 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-phenoxyphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



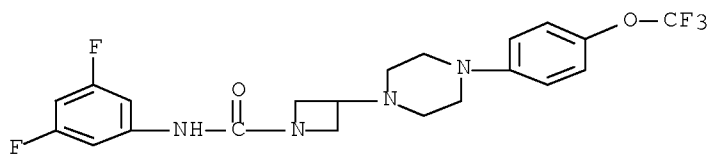
RN 858648-84-7 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-cyanophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



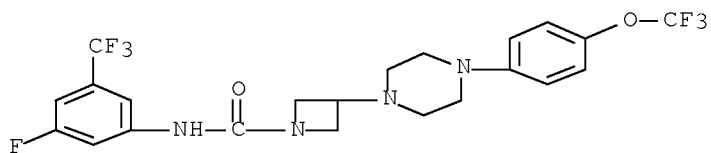
RN 858648-85-8 CAPLUS

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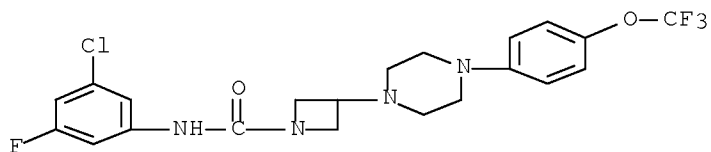
RN 858648-86-9 CAPLUS

CN 1-Azetidinecarboxamide, N-[3-fluoro-5-(trifluoromethyl)phenyl]-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)

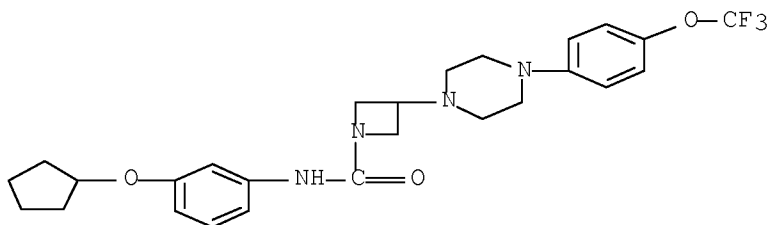


RN 858648-87-0 CAPLUS

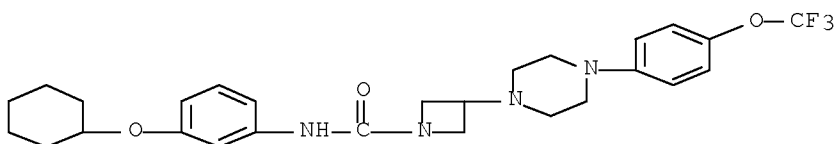
CN 1-Azetidinecarboxamide, N-(3-chloro-5-fluorophenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



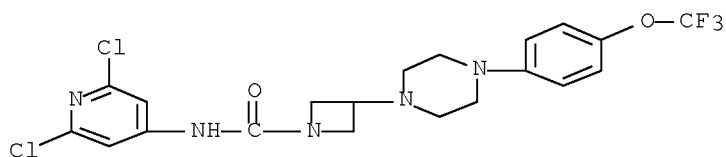
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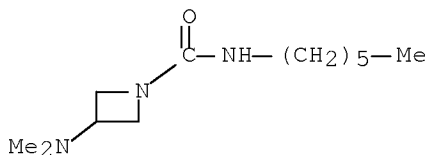
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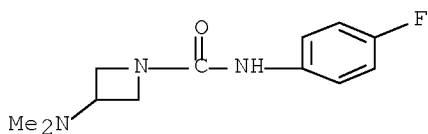
RN 858648-92-7 CAPLUS
 CN 1-Azetidinecarboxamide, N-(2,6-dichloro-4-pyridinyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



RN 858648-93-8 CAPLUS
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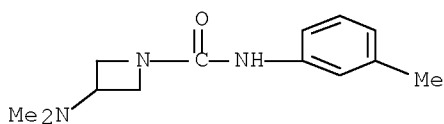


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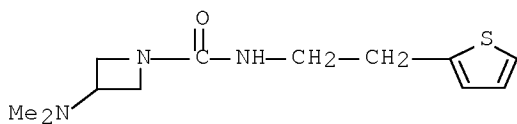
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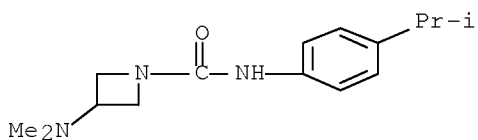
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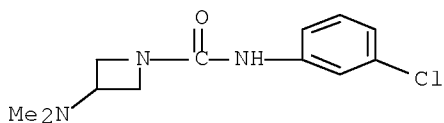
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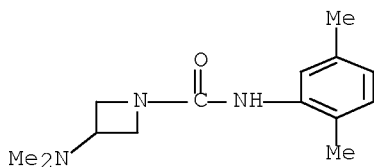
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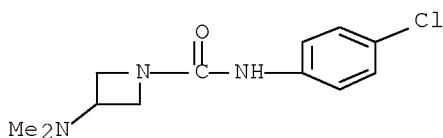
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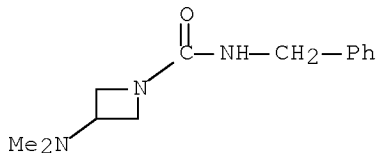
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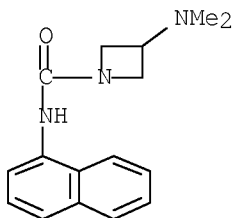
RN 858649-01-1 CAPLUS

CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(phenylmethyl)- (CA INDEX NAME)



RN 858649-02-2 CAPLUS

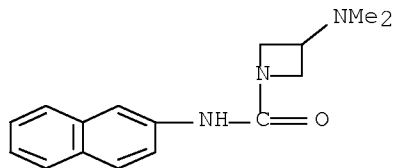
CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-1-naphthalenyl- (CA INDEX NAME)



RN 858649-03-3 CAPLUS

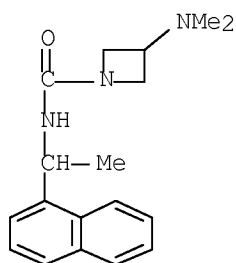
CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-2-naphthalenyl- (CA INDEX NAME)

NAME)



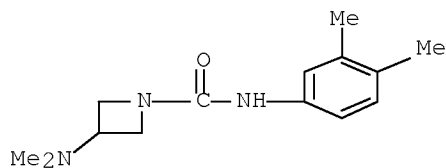
RN 858649-04-4 CAPLUS

CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-[1-(1-naphthalenyl)ethyl]-
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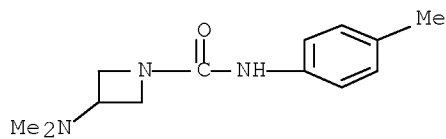
RN 858649-05-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(3,4-dimethylphenyl)- (CA
INDEX NAME)



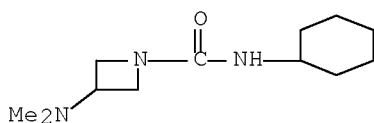
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CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(4-methylphenyl)- (CA INDEX
NAME)



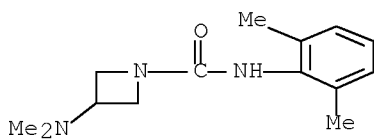
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CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(dimethylamino)- (CA INDEX NAME)



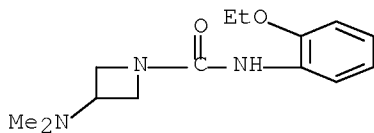
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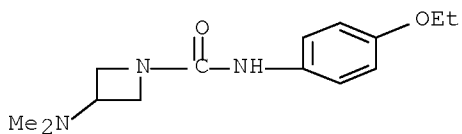
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CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(2-ethoxyphenyl)- (CA INDEX NAME)



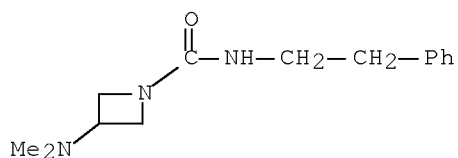
RN 858649-10-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(4-ethoxyphenyl)- (CA INDEX NAME)



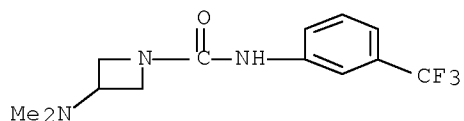
RN 858649-11-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(2-phenylethyl)- (CA INDEX NAME)



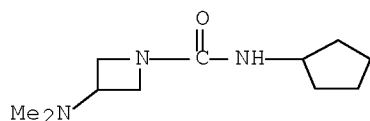
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CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-[3-(trifluoromethyl)phenyl]-
(CA INDEX NAME)



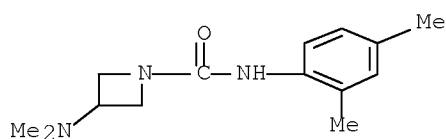
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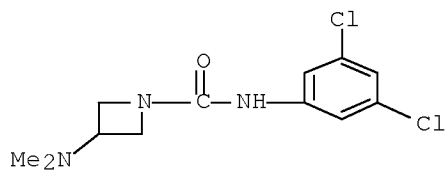
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CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(2,4-dimethylphenyl)-
(CA INDEX NAME)

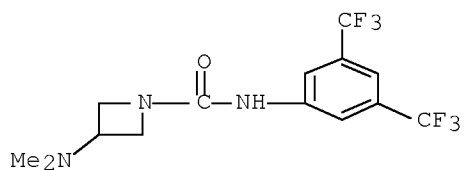


RN 858649-15-7 CAPLUS

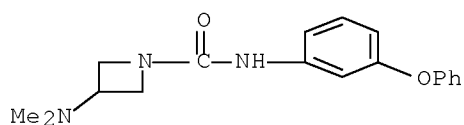
CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(dimethylamino)-
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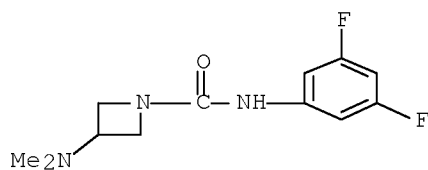
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(dimethylamino)- (CA INDEX NAME)



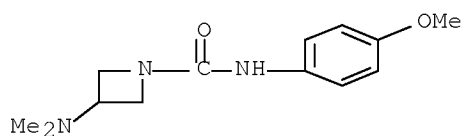
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CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(3-phenoxyphenyl)- (CA INDEX NAME)



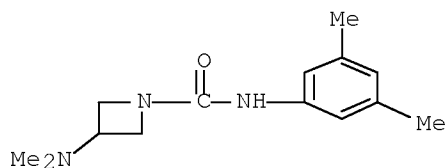
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CN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-(dimethylamino)- (CA INDEX NAME)



RN 858649-19-1 CAPLUS
CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(4-methoxyphenyl)- (CA INDEX NAME)

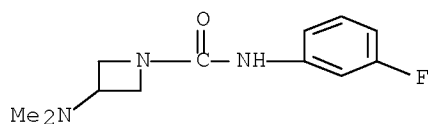


RN 858649-20-4 CAPLUS
CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(3,5-dimethylphenyl)- (CA INDEX NAME)



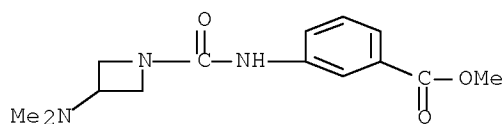
RN 858649-21-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-(dimethylamino)-N-(3-fluorophenyl)- (CA INDEX NAME)



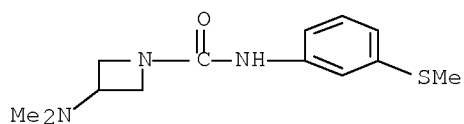
RN 858649-22-6 CAPLUS

CN Benzoic acid, 3-[[[3-(dimethylamino)-1-azetidinyl]carbonyl]amino]-, methyl ester (CA INDEX NAME)



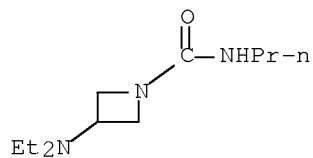
RN 858649-23-7 CAPLUS

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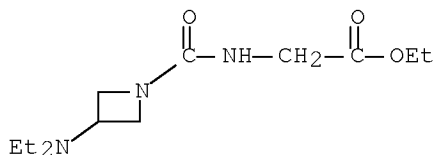
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CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-propyl- (CA INDEX NAME)



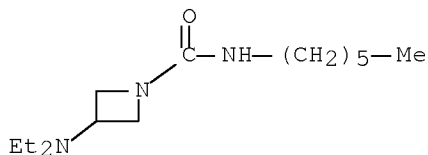
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CN Glycine, N-[[3-(diethylamino)-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)



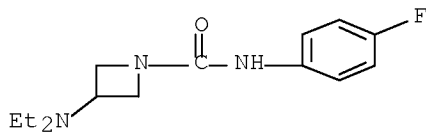
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CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-hexyl- (CA INDEX NAME)



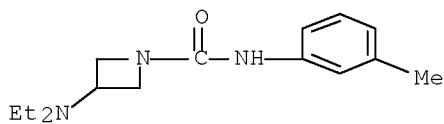
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CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(4-fluorophenyl)- (CA INDEX NAME)



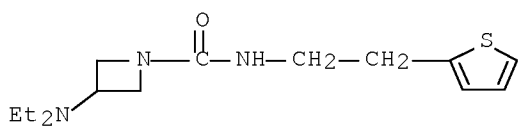
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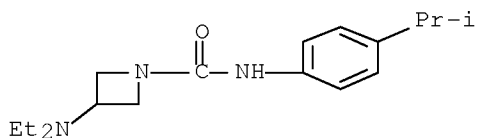
RN 858649-29-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-[2-(2-thienyl)ethyl]- (CA INDEX NAME)



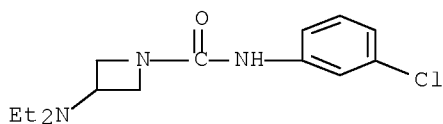
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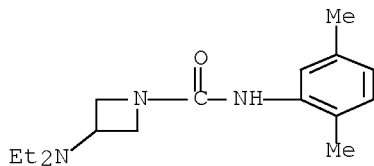
RN 858649-31-7 CAPLUS

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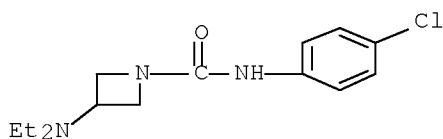
RN 858649-32-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(2,5-dimethylphenyl)- (CA INDEX NAME)



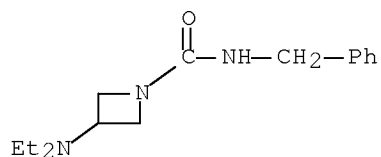
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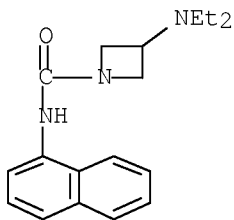
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CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(phenylmethyl)- (CA INDEX NAME)



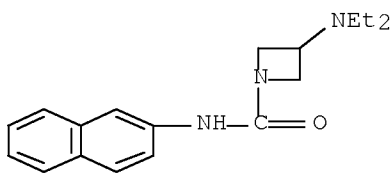
RN 858649-35-1 CAPLUS

CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-1-naphthalenyl- (CA INDEX NAME)



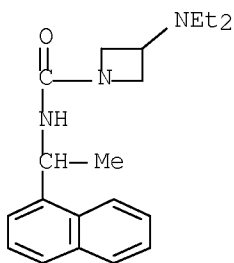
RN 858649-36-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-2-naphthalenyl- (CA INDEX NAME)

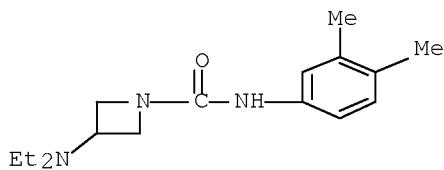


RN 858649-37-3 CAPLUS

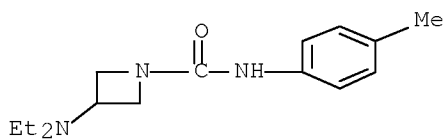
CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-[1-(1-naphthalenyl)ethyl]- (CA INDEX NAME)



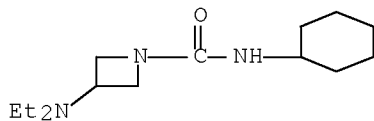
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CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(3,4-dimethylphenyl)- (CA INDEX NAME)



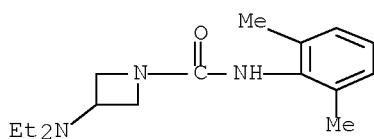
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CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(4-methylphenyl)- (CA INDEX NAME)



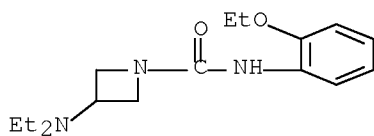
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CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(diethylamino)- (CA INDEX NAME)



RN 858649-41-9 CAPLUS
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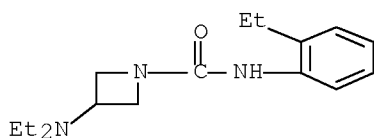


RN 858649-42-0 CAPLUS
CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(2-ethoxyphenyl)- (CA INDEX NAME)



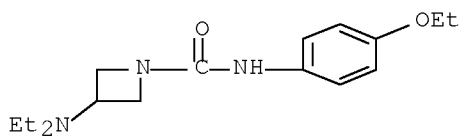
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CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(2-ethylphenyl)- (CA INDEX NAME)



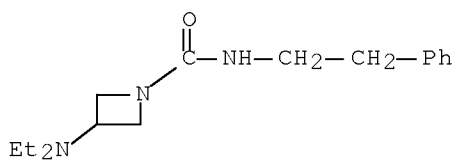
RN 858649-44-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(4-ethoxyphenyl)- (CA INDEX NAME)



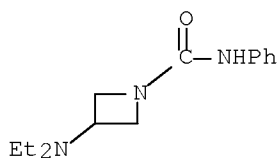
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CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(2-phenylethyl)- (CA INDEX NAME)

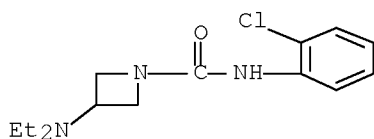


RN 858649-46-4 CAPLUS

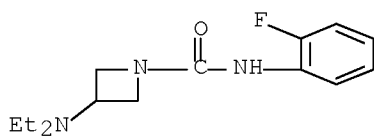
CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-phenyl- (CA INDEX NAME)



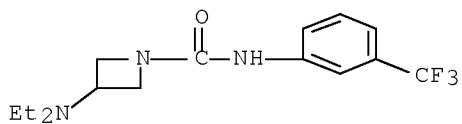
RN 858649-47-5 CAPLUS
CN 1-Azetidinecarboxamide, N-(2-chlorophenyl)-3-(diethylamino)- (CA INDEX NAME)



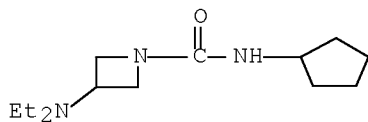
RN 858649-48-6 CAPLUS
CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(2-fluorophenyl)- (CA INDEX NAME)



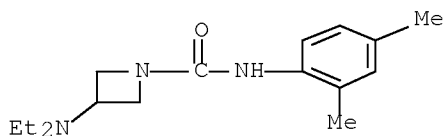
RN 858649-49-7 CAPLUS
CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



RN 858649-50-0 CAPLUS
CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(diethylamino)- (CA INDEX NAME)

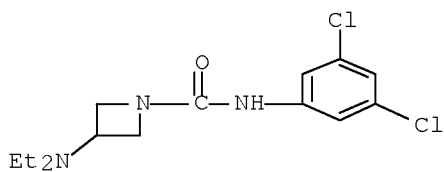


RN 858649-51-1 CAPLUS
CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(2,4-dimethylphenyl)- (CA INDEX NAME)



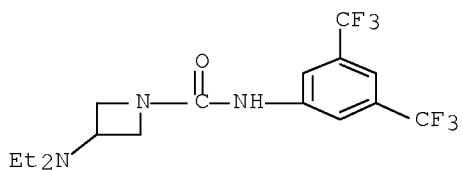
RN 858649-52-2 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(diethylamino)- (CA INDEX NAME)



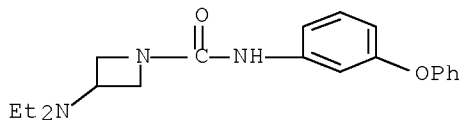
RN 858649-53-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(diethylamino)- (CA INDEX NAME)



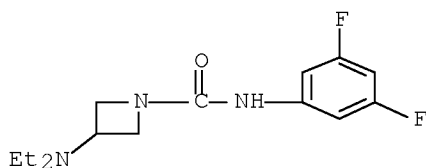
RN 858649-54-4 CAPLUS

CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(3-phenoxyphenyl)- (CA INDEX NAME)

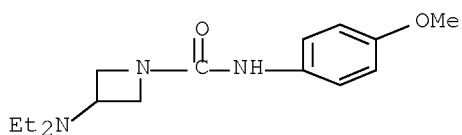


RN 858649-55-5 CAPLUS

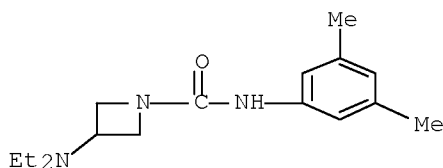
CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(3,5-difluorophenyl)- (CA INDEX NAME)



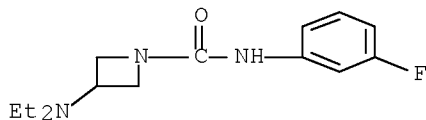
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 CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(4-methoxyphenyl)- (CA INDEX NAME)



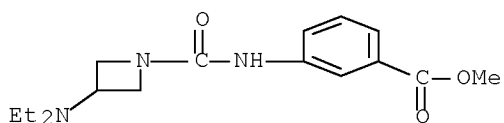
RN 858649-57-7 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(3,5-dimethylphenyl)- (CA INDEX NAME)



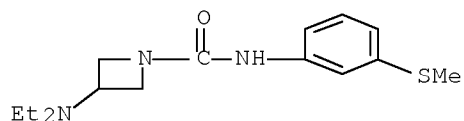
RN 858649-58-8 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(3-fluorophenyl)- (CA INDEX NAME)



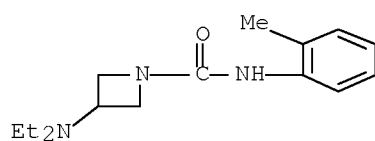
RN 858649-59-9 CAPLUS
 CN Benzoic acid, 3-[[[3-(diethylamino)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



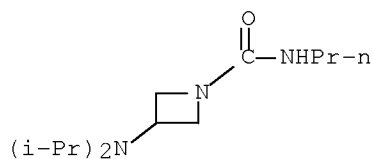
RN 858649-60-2 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-[3-(methylthio)phenyl]- (CA INDEX NAME)



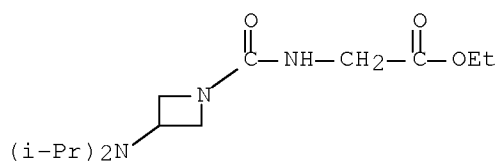
RN 858649-61-3 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(diethylamino)-N-(2-methylphenyl)- (CA INDEX NAME)



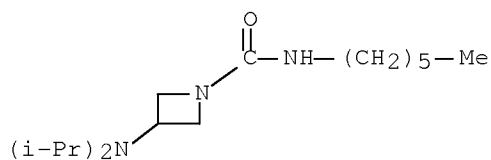
RN 858649-62-4 CAPLUS
 CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-propyl- (CA INDEX NAME)



RN 858649-63-5 CAPLUS
 CN Glycine, N-[[3-[bis(1-methylethyl)amino]-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)

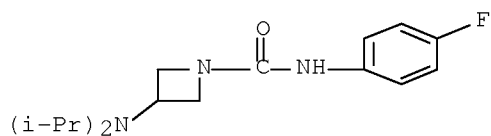


RN 858649-64-6 CAPLUS
 CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-hexyl- (CA INDEX NAME)



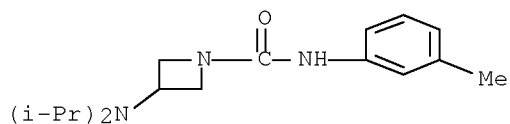
RN 858649-65-7 CAPLUS

CN 1-Azetidinecarboxamide, 3-[[bis(1-methylethyl)amino]-N-(4-fluorophenyl)]-
(CA INDEX NAME)



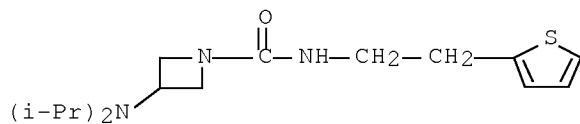
RN 858649-66-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-[[bis(1-methylethyl)amino]-N-(3-methylphenyl)]-
(CA INDEX NAME)



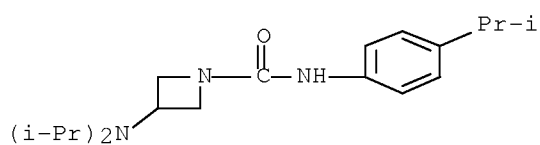
RN 858649-67-9 CAPLUS

CN 1-Azetidinecarboxamide, 3-[[bis(1-methylethyl)amino]-N-[2-(2-thienyl)ethyl]]-
(CA INDEX NAME)

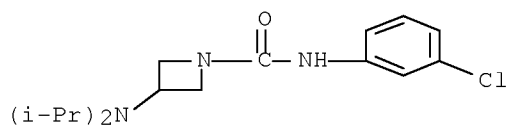


RN 858649-68-0 CAPLUS

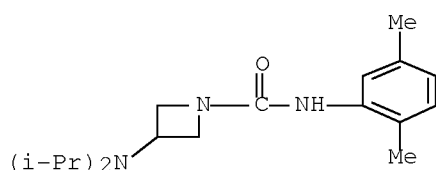
CN 1-Azetidinecarboxamide, 3-[[bis(1-methylethyl)amino]-N-[4-(1-methylethyl)phenyl]]-
(CA INDEX NAME)



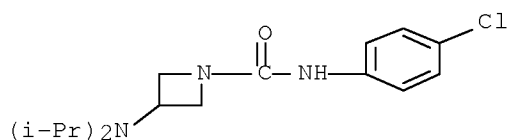
RN 858649-69-1 CAPLUS
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(3-chlorophenyl)-
(CA INDEX NAME)



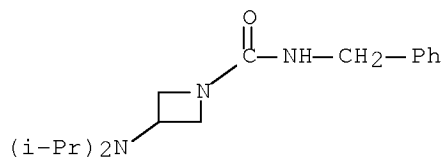
RN 858649-70-4 CAPLUS
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(2,5-dimethylphenyl)-
(CA INDEX NAME)



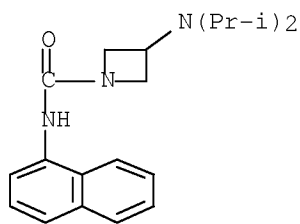
RN 858649-71-5 CAPLUS
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(4-chlorophenyl)-
(CA INDEX NAME)



RN 858649-72-6 CAPLUS
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(phenylmethyl)- (CA
INDEX NAME)

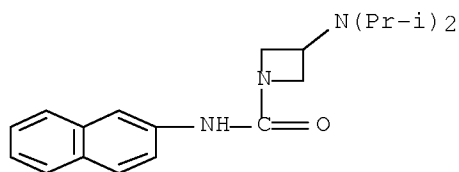


RN 858649-73-7 CAPLUS
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-1-naphthalenyl- (CA
INDEX NAME)



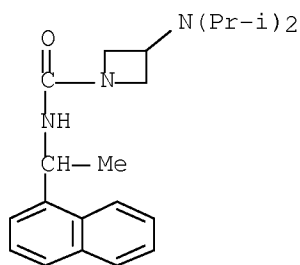
RN 858649-74-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-2-naphthalenyl- (CA INDEX NAME)



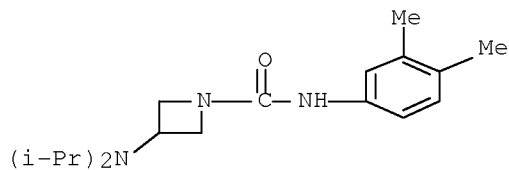
RN 858649-75-9 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-[1-(1-naphthalenyl)ethyl]- (CA INDEX NAME)



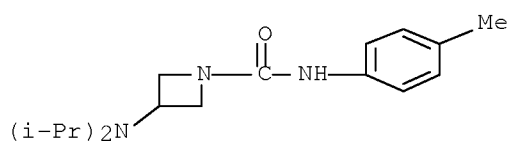
RN 858649-76-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(3,4-dimethylphenyl)- (CA INDEX NAME)



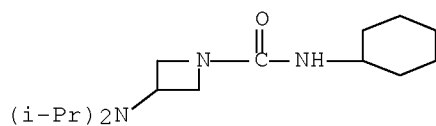
RN 858649-77-1 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(4-methylphenyl)- (CA INDEX NAME)



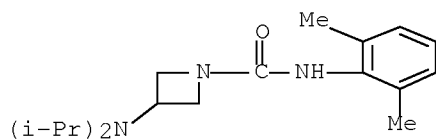
RN 858649-78-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-cyclohexyl- (CA INDEX NAME)



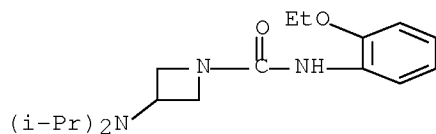
RN 858649-79-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(2,6-dimethylphenyl)- (CA INDEX NAME)



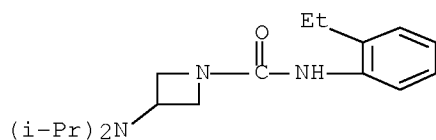
RN 858649-80-6 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(2-ethoxyphenyl)- (CA INDEX NAME)



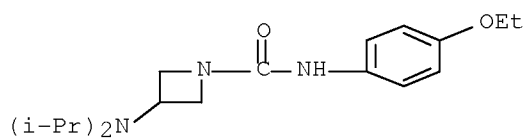
RN 858649-81-7 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(2-ethylphenyl)- (CA INDEX NAME)



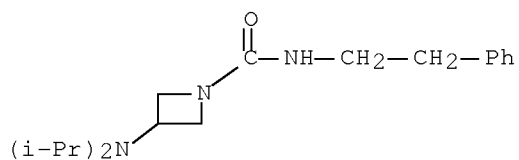
RN 858649-82-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(4-ethoxyphenyl)-
(CA INDEX NAME)



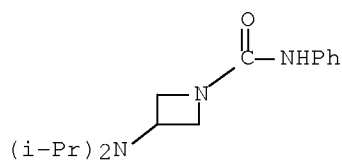
RN 858649-83-9 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(2-phenylethyl)-
(CA INDEX NAME)



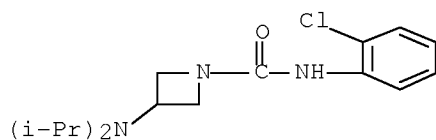
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CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-phenyl- (CA INDEX
NAME)

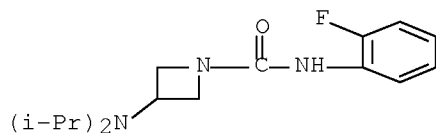


RN 858649-85-1 CAPLUS

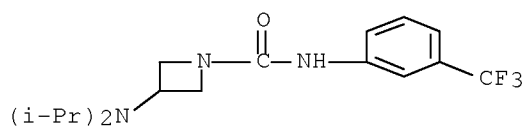
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(2-chlorophenyl)-
(CA INDEX NAME)



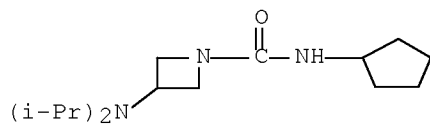
RN 858649-86-2 CAPLUS
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(2-fluorophenyl)-
(CA INDEX NAME)



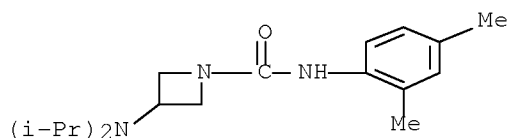
RN 858649-87-3 CAPLUS
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



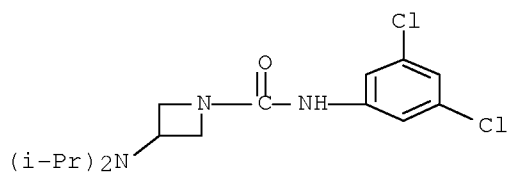
RN 858649-88-4 CAPLUS
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-cyclopentyl- (CA INDEX NAME)



RN 858649-89-5 CAPLUS
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(2,4-dimethylphenyl)-
(CA INDEX NAME)

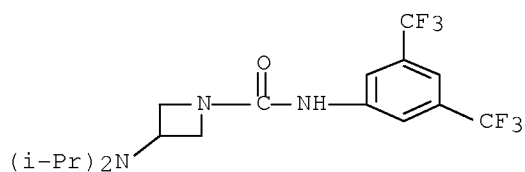


RN 858649-90-8 CAPLUS
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(3,5-dichlorophenyl)-
(CA INDEX NAME)



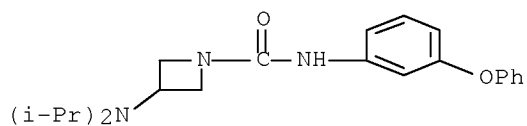
RN 858649-91-9 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-[3,5-bis(trifluoromethyl)phenyl]- (CA INDEX NAME)



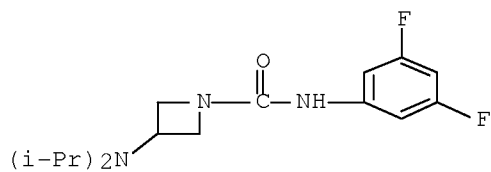
RN 858649-92-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(3-phenoxyphenyl)- (CA INDEX NAME)



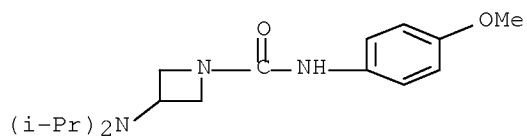
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CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(3,5-difluorophenyl)- (CA INDEX NAME)



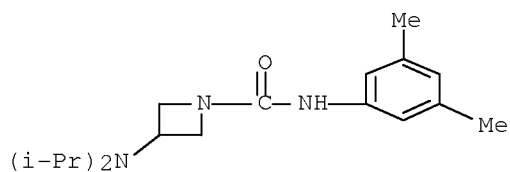
RN 858649-94-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(4-methoxyphenyl)- (CA INDEX NAME)



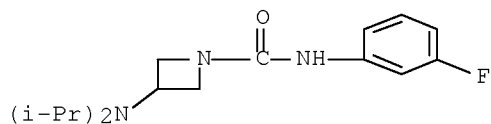
RN 858649-95-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(3,5-dimethylphenyl)-
(CA INDEX NAME)



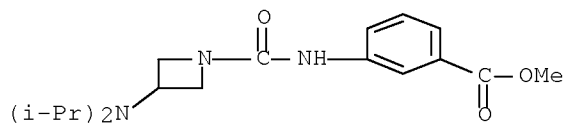
RN 858649-96-4 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(3-fluorophenyl)-
(CA INDEX NAME)



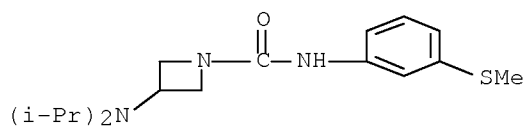
RN 858649-97-5 CAPLUS

CN Benzoic acid, 3-[[[3-[bis(1-methylethyl)amino]-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)

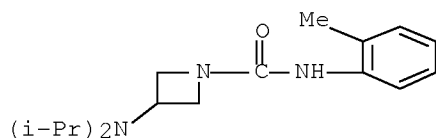


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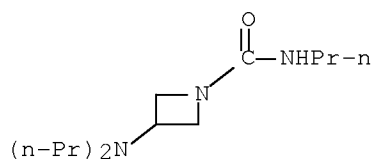
CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-[3-(methylthio)phenyl]- (CA INDEX NAME)



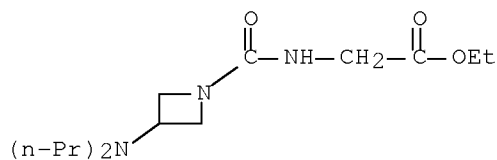
RN 858649-99-7 CAPLUS
 CN 1-Azetidinecarboxamide, 3-[bis(1-methylethyl)amino]-N-(2-methylphenyl)-
 (CA INDEX NAME)



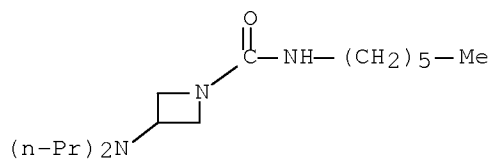
RN 858650-00-7 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-propyl- (CA INDEX NAME)



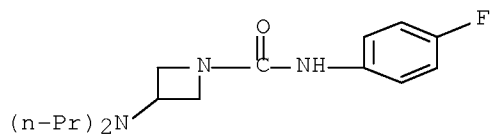
RN 858650-01-8 CAPLUS
 CN Glycine, N-[[3-(dipropylamino)-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)



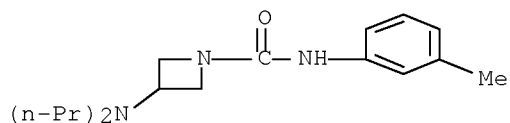
RN 858650-02-9 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-hexyl- (CA INDEX NAME)



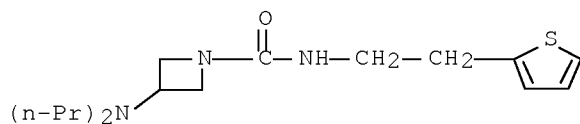
RN 858650-03-0 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-(4-fluorophenyl)- (CA INDEX NAME)



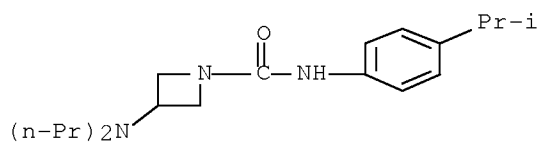
RN 858650-04-1 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-(3-methylphenyl)- (CA INDEX NAME)



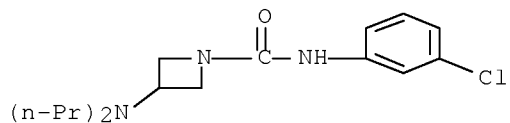
RN 858650-05-2 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-[2-(2-thienyl)ethyl]- (CA INDEX NAME)



RN 858650-06-3 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-[4-(1-methylethyl)phenyl]- (CA INDEX NAME)

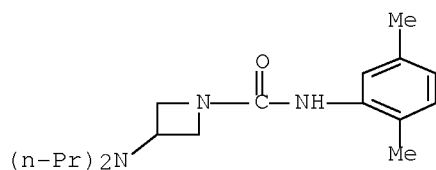


RN 858650-07-4 CAPLUS
 CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-(dipropylamino)- (CA INDEX NAME)



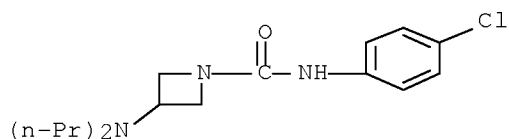
RN 858650-08-5 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,5-dimethylphenyl)-3-(dipropylamino)- (CA INDEX NAME)



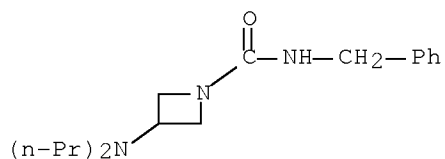
RN 858650-09-6 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-(dipropylamino)- (CA INDEX NAME)



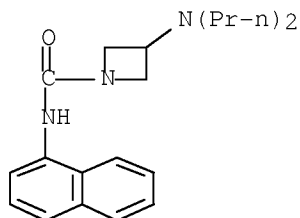
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CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-(phenylmethyl)- (CA INDEX NAME)



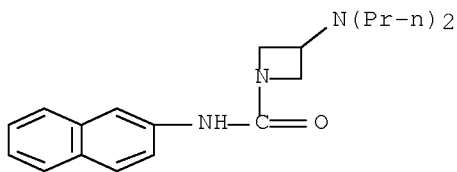
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CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-1-naphthalenyl- (CA INDEX NAME)



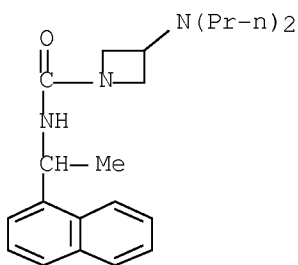
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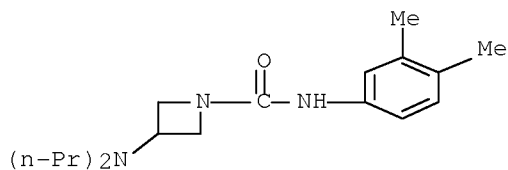
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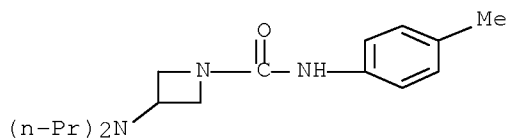
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CN 1-Azetidinecarboxamide, N-(3,4-dimethylphenyl)-3-(dipropylamino)- (CA
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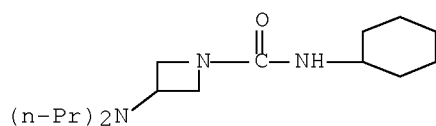
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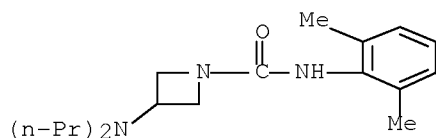
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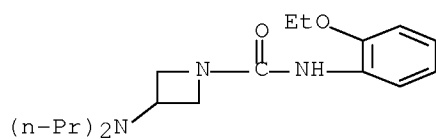
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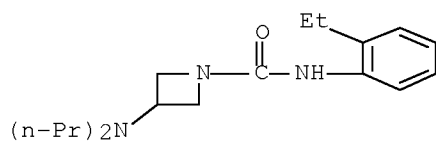
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CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-(2-ethoxyphenyl)- (CA INDEX NAME)



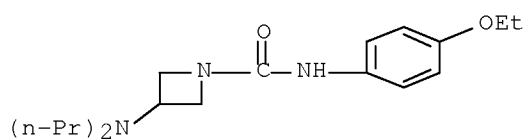
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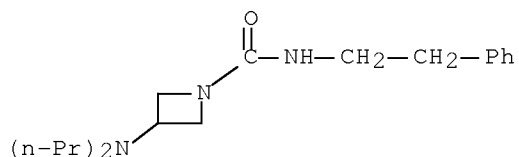


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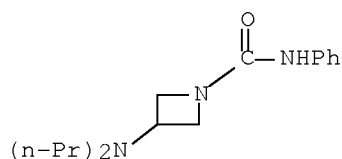
CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-(4-ethoxyphenyl)- (CA INDEX NAME)



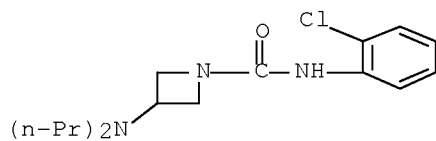
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 CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-(2-phenylethyl)- (CA INDEX NAME)



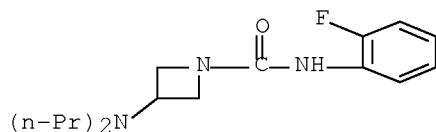
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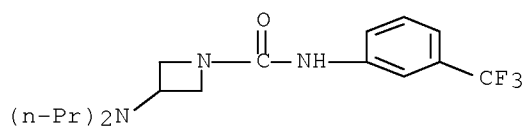
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 CN 1-Azetidinecarboxamide, N-(2-chlorophenyl)-3-(dipropylamino)- (CA INDEX NAME)



RN 858650-24-5 CAPLUS
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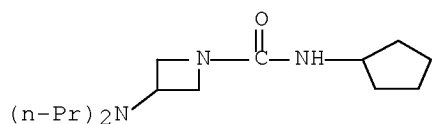


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 CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



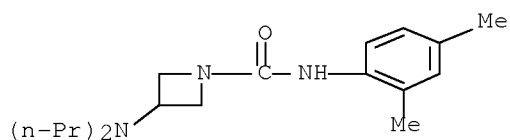
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CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(dipropylamino)- (CA INDEX NAME)



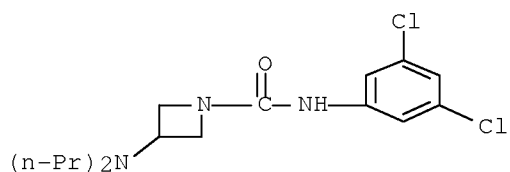
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CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(dipropylamino)- (CA INDEX NAME)



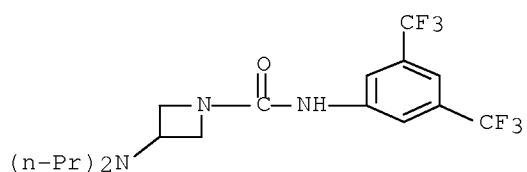
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CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(dipropylamino)- (CA INDEX NAME)

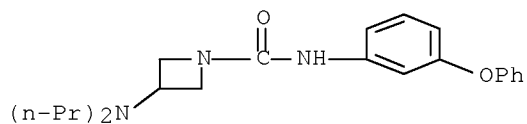


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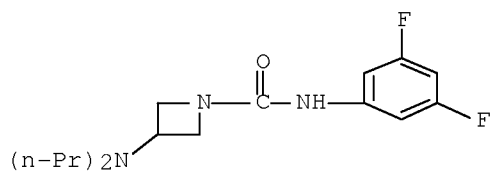
CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(dipropylamino)- (CA INDEX NAME)



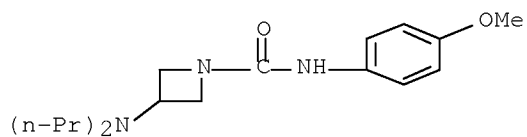
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CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-(3-phenoxyphenyl)- (CA INDEX NAME)



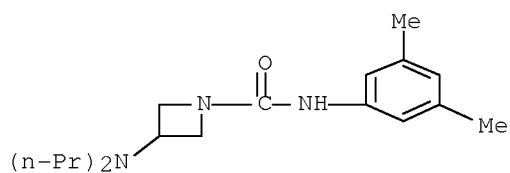
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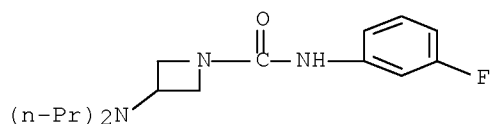
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CN 1-Azetidinecarboxamide, 3-(dipropylamino)-N-(4-methoxyphenyl)- (CA INDEX NAME)



RN 858650-36-9 CAPLUS
CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-(dipropylamino)- (CA INDEX NAME)

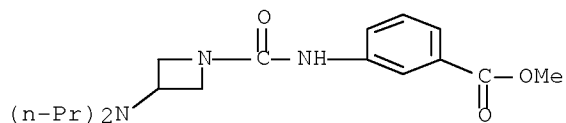


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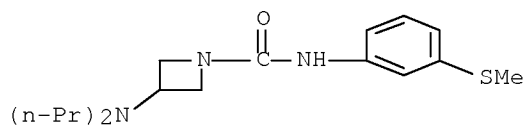
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CN Benzoic acid, 3-[[[3-(dipropylamino)-1-azetidinyl]carbonyl]amino]-, methyl ester (CA INDEX NAME)



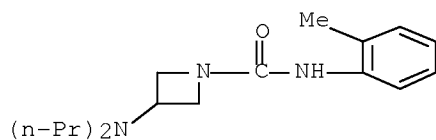
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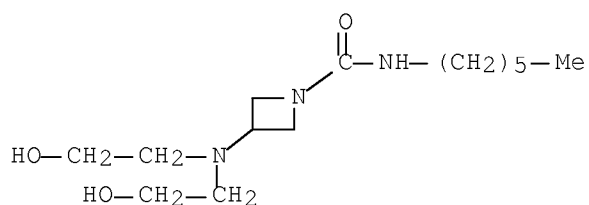
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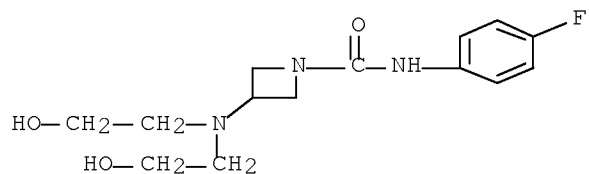
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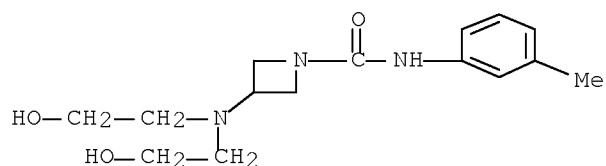
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(CA INDEX NAME)



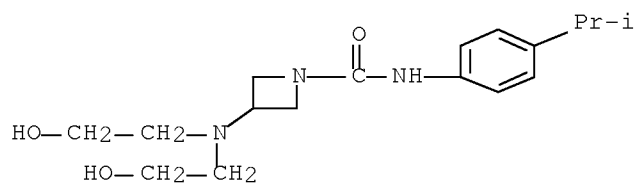
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-(3-methylphenyl)-
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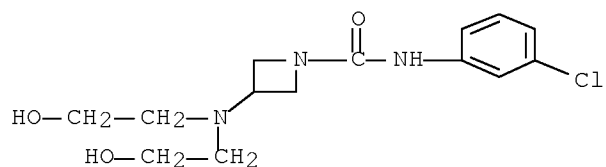
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-[4-(1-methylethyl)phenyl]- (CA INDEX NAME)



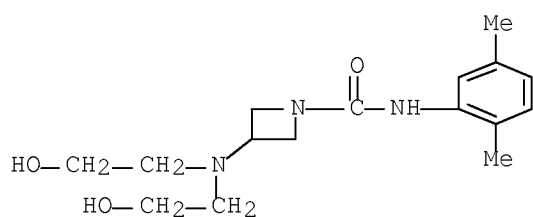
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-(3-chlorophenyl)-
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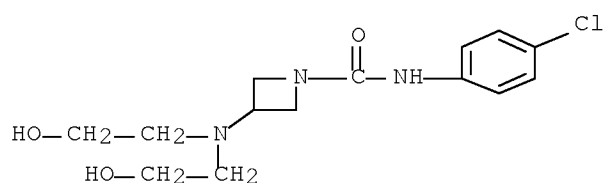
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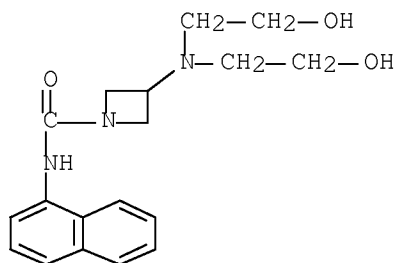
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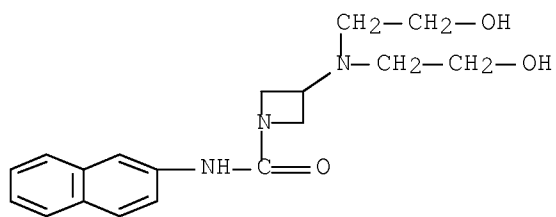
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-1-naphthalenyl- (CA INDEX NAME)



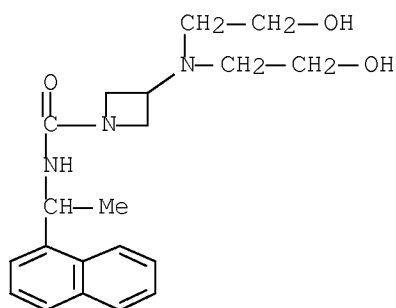
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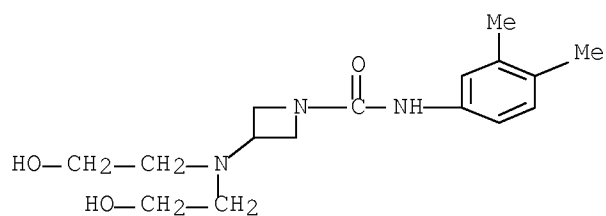
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-[1-(1-naphthalenyl)ethyl]- (CA INDEX NAME)



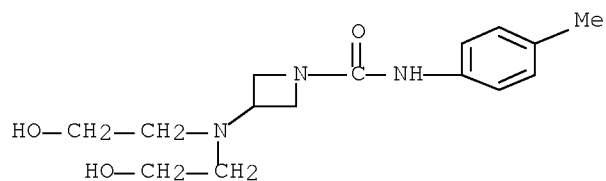
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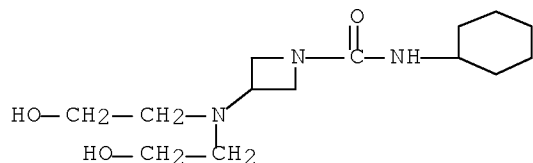
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-(4-methylphenyl)- (CA INDEX NAME)



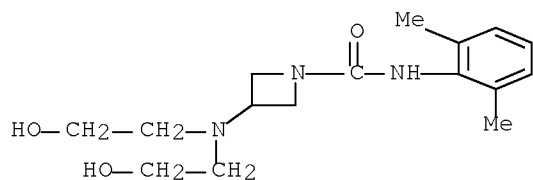
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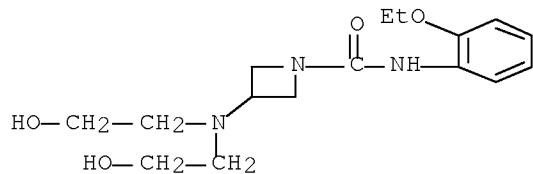
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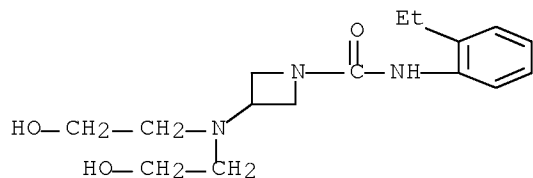
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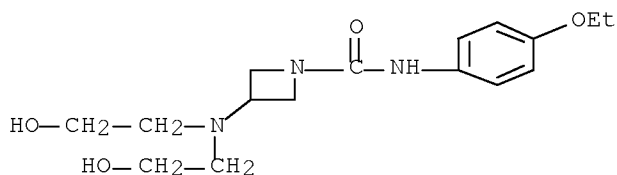
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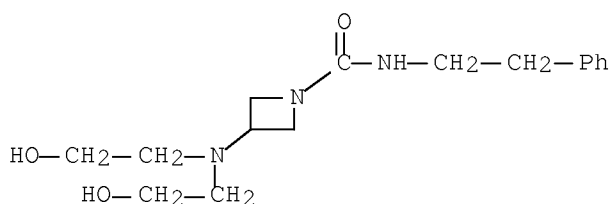
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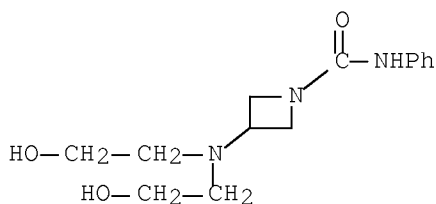
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-(2-phenylethyl)-
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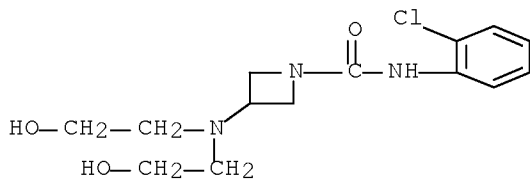
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-phenyl- (CA INDEX
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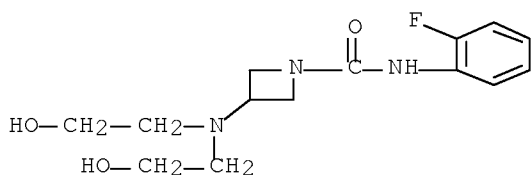
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-(2-chlorophenyl)-
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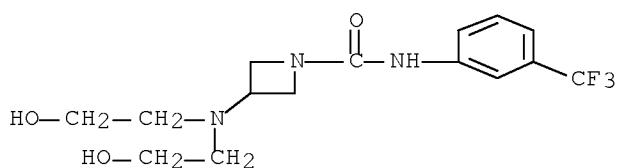
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-(2-fluorophenyl)-
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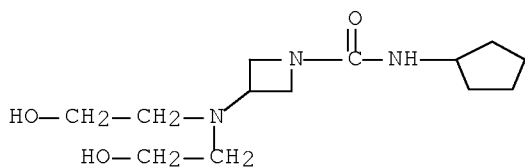
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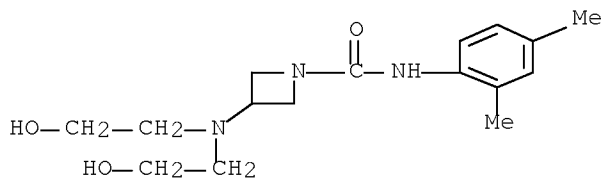
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-cyclopentyl- (CA INDEX NAME)



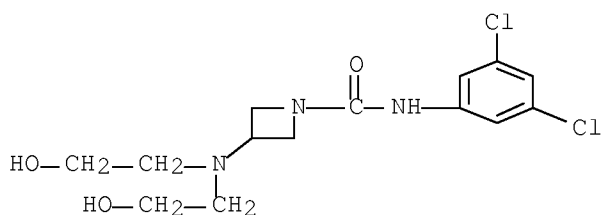
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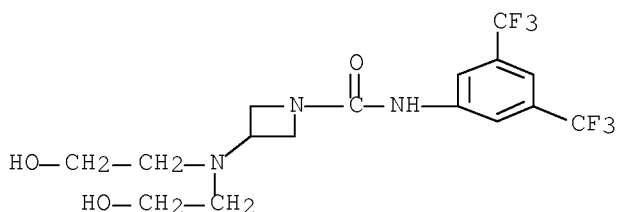
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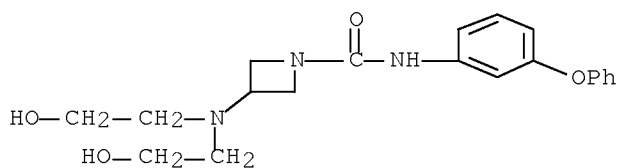
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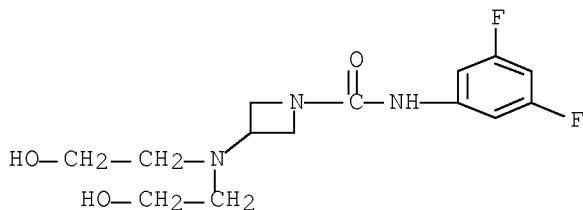
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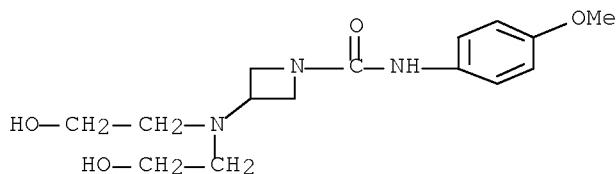
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-(3,5-difluorophenyl)- (CA INDEX NAME)



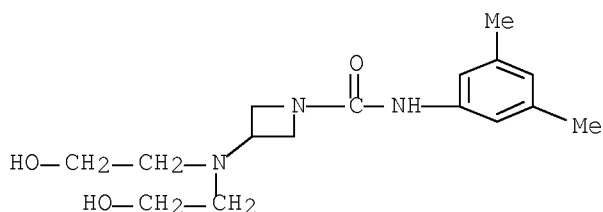
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CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-(4-methoxyphenyl)- (CA INDEX NAME)



RN 858650-77-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-(3,5-dimethylphenyl)- (CA INDEX NAME)



IT 858650-78-9P, 3-[Bis(2-hydroxyethyl)amino]-N-(3-fluorophenyl)azetidine-1-carboxamide 858650-79-0P, Methyl 3-[[[3-[bis(2-hydroxyethyl)amino]azetidin-1-yl]carbonyl]amino]benzoate 858650-80-3P, 3-[Bis(2-hydroxyethyl)amino]-N-[3-(methylsulfanyl)phenyl]azetidine-1-carboxamide 858650-81-4P, N-Ethyl-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858650-82-5P, 3-[Methyl(phenyl)amino]-N-propylazetidine-1-carboxamide 858650-83-6P 858650-84-7P, N-Hexyl-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858650-85-8P, N-(4-Fluorophenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858650-86-9P, N-(3-Methylphenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858650-88-1P, 3-[Methyl(phenyl)amino]-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858650-90-5P, N-(4-Isopropylphenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858650-91-6P, N-(3-Chlorophenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858650-92-7P, N-(2,5-Dimethylphenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858650-94-9P, N-(4-Chlorophenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858650-96-1P, N-Benzyl-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858650-98-3P, 3-[Methyl(phenyl)amino]-N-(1-naphthyl)azetidine-1-carboxamide 858650-99-4P, 3-[Methyl(phenyl)amino]-N-(2-naphthyl)azetidine-1-carboxamide 858651-00-0P, 3-[Methyl(phenyl)amino]-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858651-01-1P, N-(3,4-Dimethylphenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858651-02-2P, N-(4-Methylphenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858651-03-3P, N-Cyclohexyl-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858651-04-4P, N-(2,6-Dimethylphenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858651-06-6P, N-(2-Ethoxyphenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858651-08-8P, N-(2-Ethylphenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide 858651-10-2P, N-(4-Ethoxyphenyl)-3-[methyl(phenyl)amino]azetidine-1-carboxamide

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858651-77-1P, 3-[Ethyl(phenyl)amino]-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858651-78-2P, N-(3,5-Difluorophenyl)-3-[ethyl(phenyl)amino]azetidine-1-carboxamide 858651-79-3P, 3-[Ethyl(phenyl)amino]-N-(4-methoxyphenyl)azetidine-1-carboxamide 858651-80-6P, N-(3,5-Dimethylphenyl)-3-[ethyl(phenyl)amino]azetidine-1-carboxamide 858651-81-7P, 3-[Ethyl(phenyl)amino]-N-(3-fluorophenyl)azetidine-1-carboxamide 858651-82-8P, Methyl 3-[[[3-[ethyl(phenyl)amino]azetidin-1-yl]carbonyl]amino]benzoate 858651-83-9P, 3-[Ethyl(phenyl)amino]-N-[3-(methylsulfonyl)phenyl]azetidine-1-carboxamide 858651-84-0P, 3-[Ethyl(phenyl)amino]-N-(2-methylphenyl)azetidine-1-carboxamide 858651-85-1P, 3-[(Phenyl)(propyl)amino]-N-propylazetidine-1-carboxamide 858651-86-2P 858651-87-3P, N-Hexyl-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-88-4P, N-(4-Fluorophenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-89-5P, N-(3-Methylphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-90-8P, 3-[(Phenyl)(propyl)amino]-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858651-91-9P, N-(4-Isopropylphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-92-0P, N-(3-Chlorophenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-93-1P, N-(2,5-Dimethylphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-94-2P, N-(4-Chlorophenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-95-3P, N-Benzyl-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-96-4P, N-(1-Naphthyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-97-5P, N-(2-Naphthyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-98-6P, N-[1-(1-Naphthyl)ethyl]-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858651-99-7P, N-(3,4-Dimethylphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-00-3P, N-(4-Methylphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-01-4P, N-Cyclohexyl-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-02-5P, N-(2,6-Dimethylphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-03-6P, N-(2-Ethoxyphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-04-7P, N-(2-Ethylphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-05-8P, N-(4-Ethoxyphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-06-9P, N-(2-Phenylethyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-07-0P, N-Phenyl-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-08-1P, N-(2-Chlorophenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-09-2P, N-(2-Fluorophenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-10-5P, 3-[(Phenyl)(propyl)amino]-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858652-11-6P, N-Cyclopentyl-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-12-7P, N-(2,4-Dimethylphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-13-8P, N-(3,5-Dichlorophenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-14-9P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-15-0P, N-(3-Phenoxyphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-16-1P, N-(3,5-Difluorophenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-17-2P, N-(4-Methoxyphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-18-3P, N-(3,5-Dimethylphenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-19-4P, N-(3-Fluorophenyl)-3-[phenyl(propyl)amino]azetidine-1-carboxamide 858652-20-7P, Methyl 3-[[[3-[phenyl(propyl)amino]azetidin-1-

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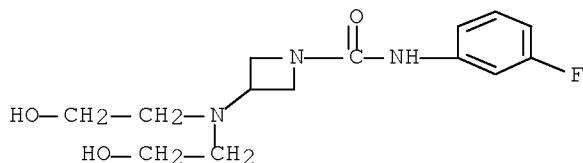
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858652-67-2P, N-(3-Methylphenyl)-3-pyrrolidin-1-ylazetidine-1-
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N-(4-Isopropylphenyl)-3-pyrrolidin-1-ylazetidine-1-carboxamide
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N-(2-Fluorophenyl)-3-pyrrolidin-1-ylazetidine-1-carboxamide
858652-88-7P, 3-Pyrrolidin-1-yl-N-[3-(trifluoromethyl)phenyl]azeti
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ylazetidine-1-carboxamide 858652-90-1P, N-(3,5-Dichlorophenyl)-3-
pyrrolidin-1-ylazetidine-1-carboxamide 858652-91-2P,
N-[3,5-Bis(trifluoromethyl)phenyl]-3-pyrrolidin-1-ylazetidine-1-
carboxamide 858652-92-3P, N-(3-Phenoxyphenyl)-3-pyrrolidin-1-
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N-(4-Methoxyphenyl)-3-pyrrolidin-1-ylazetidine-1-carboxamide
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carboxamide 858652-96-7P, N-(3-Fluorophenyl)-3-pyrrolidin-1-
ylazetidine-1-carboxamide 858652-97-8P, Methyl
3-[[3-(3-pyrrolidin-1-ylazetidin-1-yl)carbonyl]amino]benzoate
858652-98-9P, N-[3-(Methylsulfonyl)phenyl]-3-pyrrolidin-1-
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propylazetidine-1-carboxamide 858653-00-6P 858653-01-7P
, N-Hexyl-3-piperidin-1-ylazetidine-1-carboxamide 858653-02-8P,
N-(4-Fluorophenyl)-3-piperidin-1-ylazetidine-1-carboxamide
858653-03-9P, N-(3-Methylphenyl)-3-piperidin-1-ylazetidine-1-
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N-Benzyl-3-piperidin-1-ylazetidine-1-carboxamide 858653-10-8P,
N-(1-Naphthyl)-3-piperidin-1-ylazetidine-1-carboxamide
858653-11-9P, N-(2-Naphthyl)-3-piperidin-1-ylazetidine-1-
carboxamide 858653-12-0P, N-[1-(1-Naphthyl)ethyl]-3-piperidin-1-
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piperidin-1-ylazetidine-1-carboxamide 858653-14-2P,
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 ylazetidine-1-carboxamide 858653-18-6P, N-(2-Ethylphenyl)-3-
 piperidin-1-ylazetidine-1-carboxamide 858653-19-7P,
 N-(4-Ethoxyphenyl)-3-piperidin-1-ylazetidine-1-carboxamide
 858653-20-0P, N-(2-Phenylethyl)-3-piperidin-1-ylazetidine-1-
 carboxamide 858653-21-1P, N-Phenyl-3-piperidin-1-ylazetidine-1-
 carboxamide 858653-22-2P, N-(2-Chlorophenyl)-3-piperidin-1-
 ylazetidine-1-carboxamide 858653-23-3P, N-(2-Fluorophenyl)-3-
 piperidin-1-ylazetidine-1-carboxamide 858653-24-4P,
 3-Piperidin-1-yl-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide
 858653-25-5P, N-Cyclopentyl-3-piperidin-1-ylazetidine-1-
 carboxamide 858653-26-6P, N-(2,4-Dimethylphenyl)-3-piperidin-1-
 ylazetidine-1-carboxamide 858653-27-7P, N-(3,5-Dichlorophenyl)-3-
 piperidin-1-ylazetidine-1-carboxamide 858653-28-8P,
 N-[3,5-Bis(trifluoromethyl)phenyl]-3-piperidin-1-ylazetidine-1-carboxamide
 858653-29-9P, N-(3-Phenoxyphenyl)-3-piperidin-1-ylazetidine-1-
 carboxamide 858653-30-2P, N-(3,5-Difluorophenyl)-3-piperidin-1-
 ylazetidine-1-carboxamide
 858653-31-3P, N-(4-Methoxyphenyl)-3-piperidin-1-ylazetidine-1-
 carboxamide 858653-32-4P, N-(3,5-Dimethylphenyl)-3-piperidin-1-
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 piperidin-1-ylazetidine-1-carboxamide 858653-34-6P, Methyl
 3-[[[3-piperidin-1-ylazetidin-1-yl]carbonyl]amino]benzoate
 858653-35-7P, N-[3-(Methylsulfonyl)phenyl]-3-piperidin-1-
 ylazetidine-1-carboxamide 858653-36-8P, N-(2-Methylphenyl)-3-
 piperidin-1-ylazetidine-1-carboxamide 858653-37-9P,
 3-(Azepan-1-yl)-N-propylazetidine-1-carboxamide 858653-38-0P
 858653-39-1P, 3-(Azepan-1-yl)-N-hexylazetidine-1-carboxamide
 858653-40-4P, 3-(Azepan-1-yl)-N-(4-fluorophenyl)azetidine-1-
 carboxamide 858653-41-5P, 3-(Azepan-1-yl)-N-(3-
 methylphenyl)azetidine-1-carboxamide 858653-42-6P,
 3-(Azepan-1-yl)-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide
 RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
 (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
 (Uses)

(preparation of azetidine ring compds. as EDG-5 antagonists for treatment
 of hypertension, myocardial infarction, etc.)

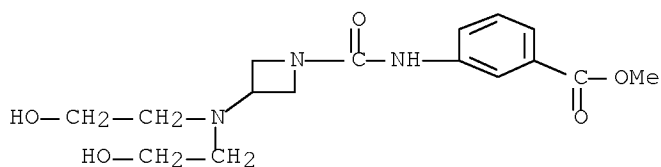
RN 858650-78-9 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(2-hydroxyethyl)amino]-N-(3-fluorophenyl)-
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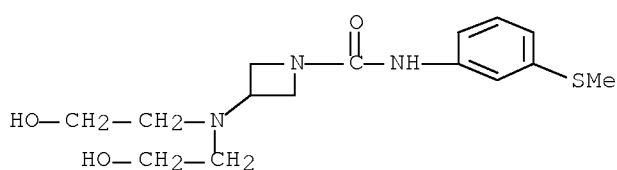
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CN Benzoic acid, 3-[[[3-[bis(2-hydroxyethyl)amino]-1-
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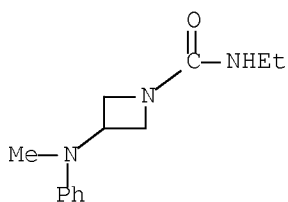
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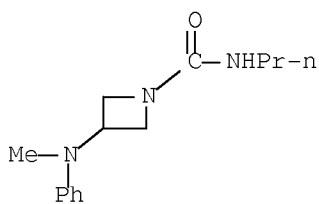
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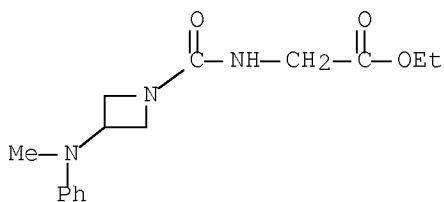
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CN 1-Azetidinecarboxamide, 3-(methylphenylamino)-N-propyl- (CA INDEX NAME)



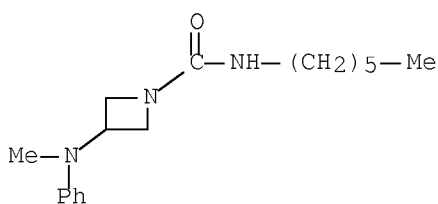
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CN Glycine, N-[[3-(methylphenylamino)-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)



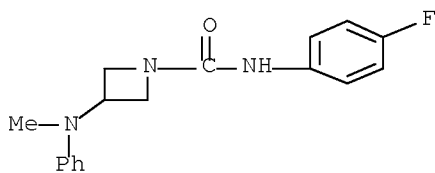
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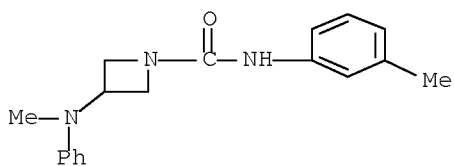
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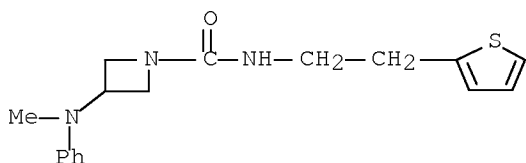
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CN 1-Azetidinecarboxamide, N-(3-methylphenyl)-3-(methylphenylamino)- (CA INDEX NAME)



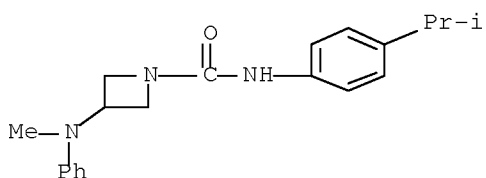
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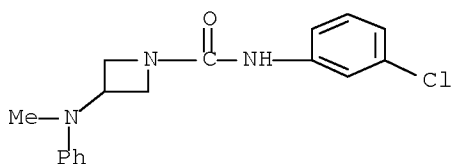
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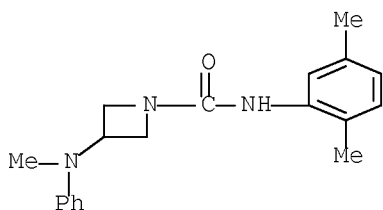
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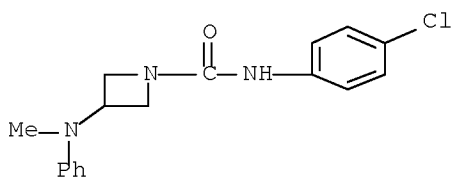
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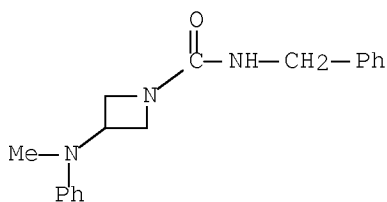


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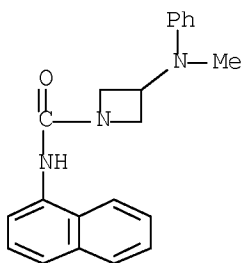
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INDEX NAME)



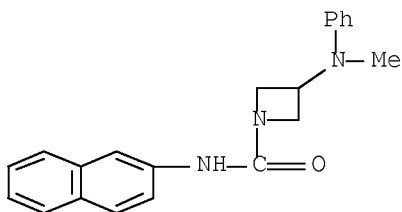
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 CN 1-Azetidinecarboxamide, 3-(methylphenylamino)-N-(phenylmethyl)- (CA INDEX NAME)



RN 858650-98-3 CAPLUS
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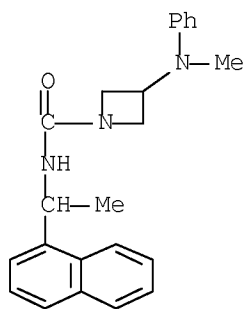


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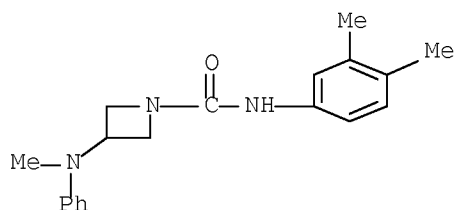
RN 858651-00-0 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(methylphenylamino)-N-[1-(1-naphthalenyl)ethyl]-

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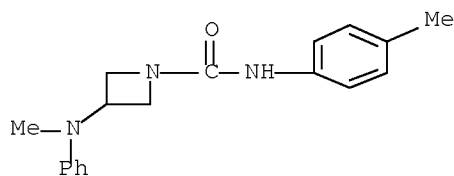
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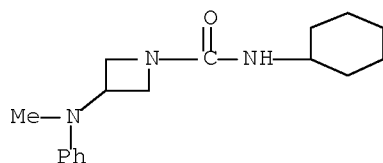
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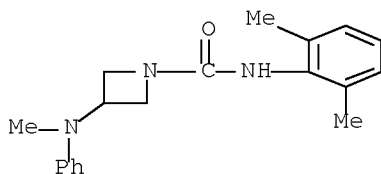
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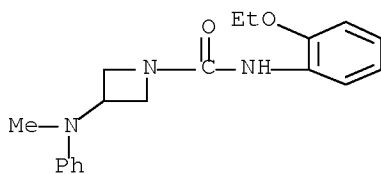
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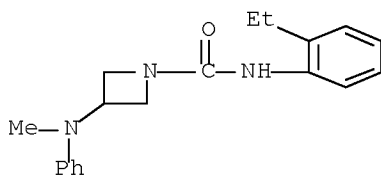
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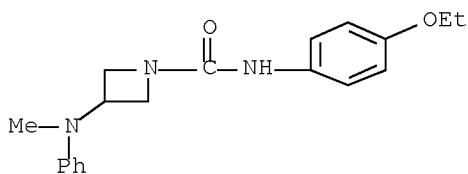
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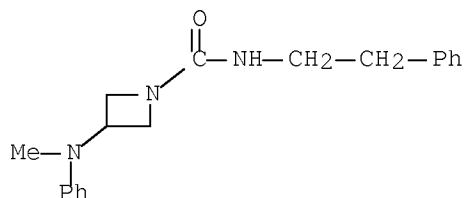
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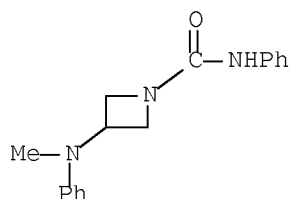
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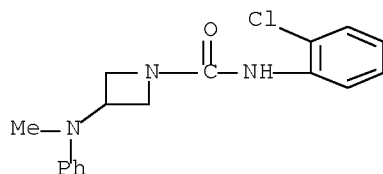
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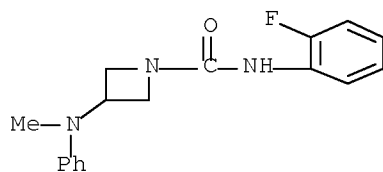
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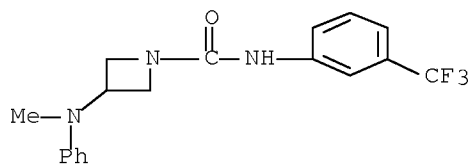
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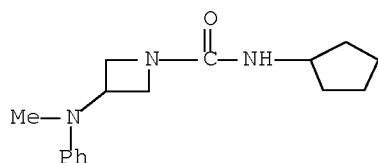
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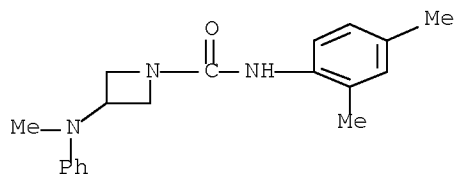
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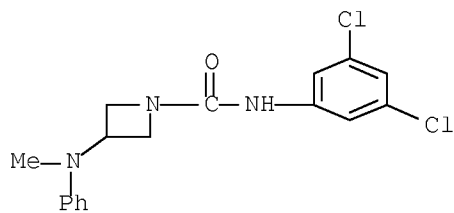
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CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(methylphenylamino)- (CA INDEX NAME)



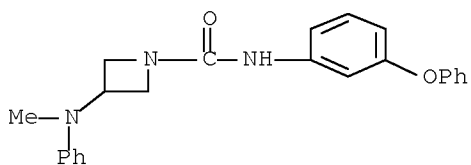
RN 858651-20-4 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(methylphenylamino)- (CA INDEX NAME)



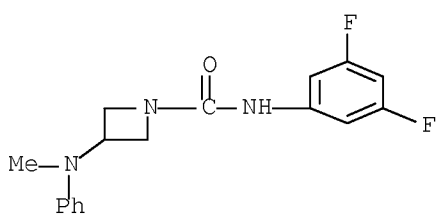
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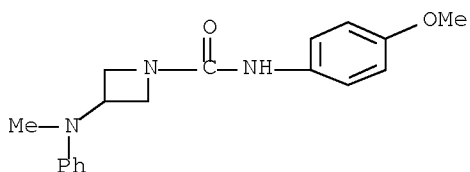
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CN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-(methylphenylamino)- (CA INDEX NAME)



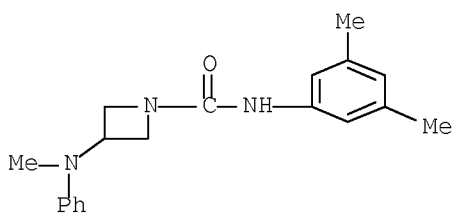
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CN 1-Azetidinecarboxamide, N-(4-methoxyphenyl)-3-(methylphenylamino)- (CA INDEX NAME)



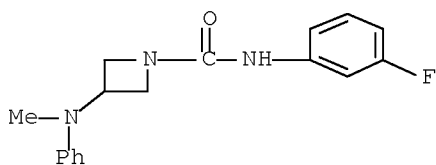
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CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-(methylphenylamino)- (CA INDEX NAME)



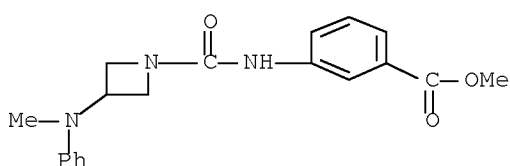
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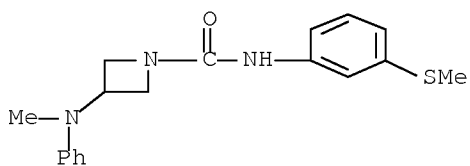
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CN Benzoic acid, 3-[[[3-(methylphenylamino)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



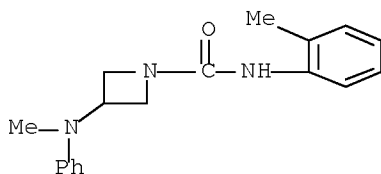
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CN 1-Azetidinecarboxamide, 3-(methylphenylamino)-N-[3-(methylthio)phenyl]- (CA INDEX NAME)



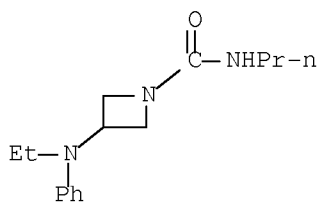
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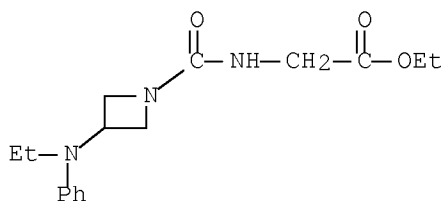


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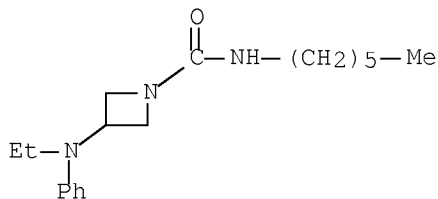
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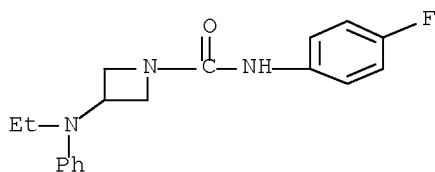
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 CN Glycine, N-[[3-(ethylphenylamino)-1-azetidiny]carbonyl]-, ethyl ester
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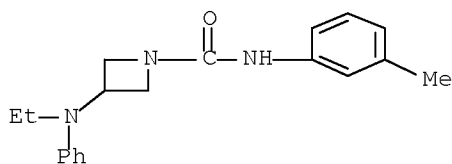
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 CN 1-Azetidinecarboxamide, 3-(ethylphenylamino)-N-hexyl- (CA INDEX NAME)



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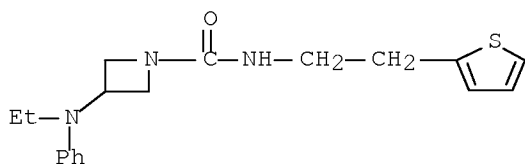


RN 858651-36-2 CAPLUS
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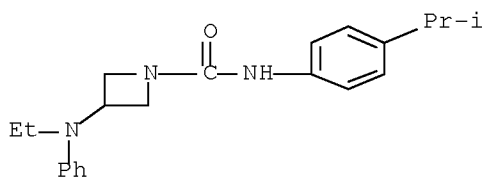
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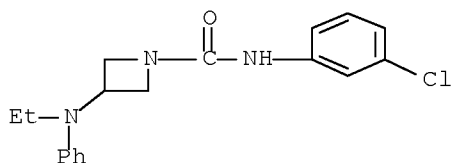
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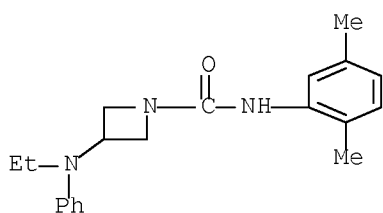
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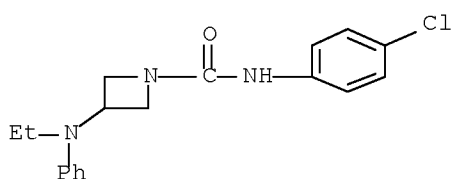
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CN 1-Azetidinecarboxamide, N-(2,5-dimethylphenyl)-3-(ethylphenylamino)- (CA INDEX NAME)



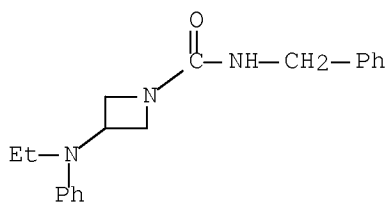
RN 858651-44-2 CAPLUS

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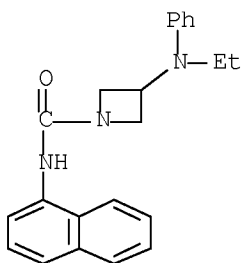
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RN 858651-47-5 CAPLUS

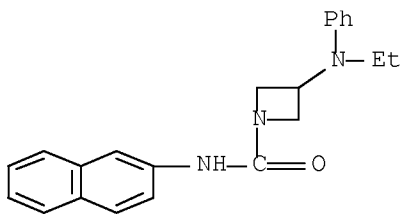
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RN 858651-49-7 CAPLUS

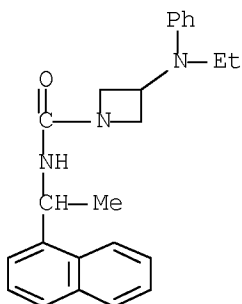
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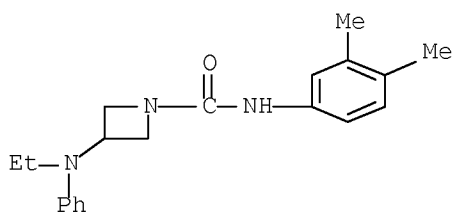
RN 858651-51-1 CAPLUS

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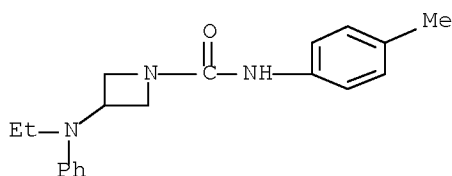
RN 858651-53-3 CAPLUS

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INDEX NAME)



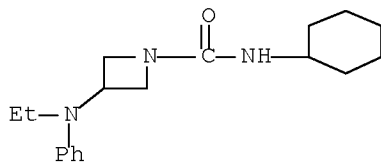
RN 858651-55-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-(ethylphenylamino)-N-(4-methylphenyl)- (CA
INDEX NAME)



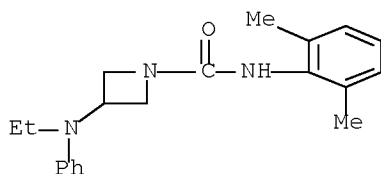
RN 858651-57-7 CAPLUS

CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(ethylphenylamino)- (CA INDEX NAME)



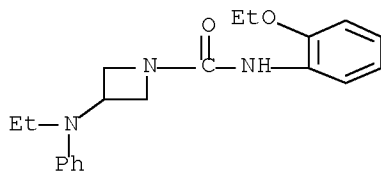
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CN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-(ethylphenylamino)- (CA INDEX NAME)



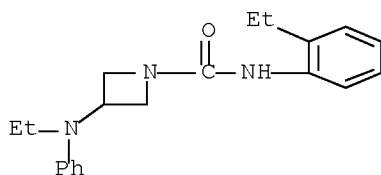
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CN 1-Azetidinecarboxamide, N-(2-ethoxyphenyl)-3-(ethylphenylamino)- (CA INDEX NAME)

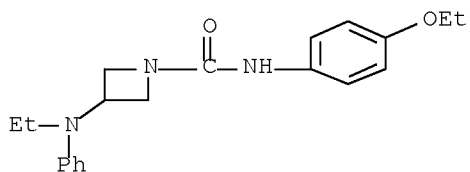


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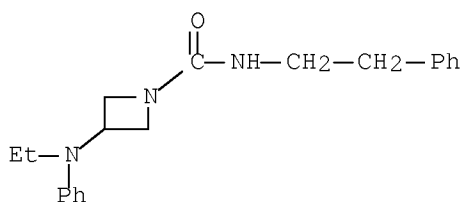
CN 1-Azetidinecarboxamide, N-(2-ethylphenyl)-3-(ethylphenylamino)- (CA INDEX NAME)



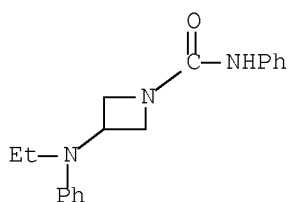
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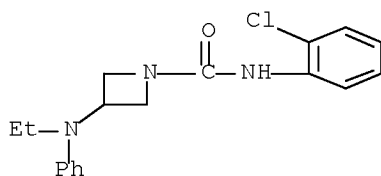
RN 858651-65-7 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(ethylphenylamino)-N-(2-phenylethyl)- (CA INDEX NAME)



RN 858651-66-8 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(ethylphenylamino)-N-phenyl- (CA INDEX NAME)

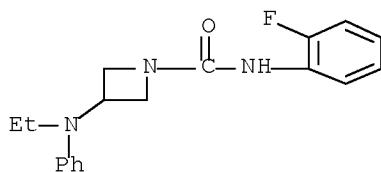


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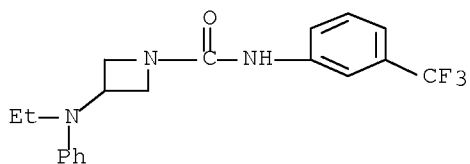
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CN 1-Azetidinecarboxamide, 3-(ethylphenylamino)-N-(2-fluorophenyl)- (CA
INDEX NAME)



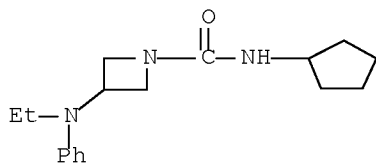
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CN 1-Azetidinecarboxamide, 3-(ethylphenylamino)-N-[3-(trifluoromethyl)phenyl]-
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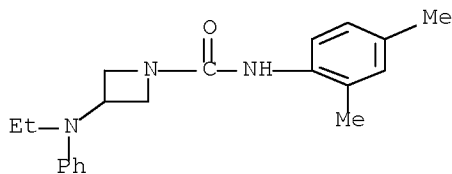
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CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(ethylphenylamino)- (CA INDEX
NAME)



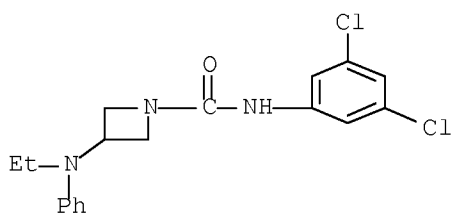
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CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(ethylphenylamino)- (CA
INDEX NAME)



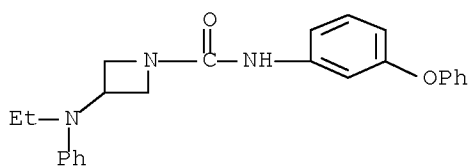
RN 858651-75-9 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(ethylphenylamino)- (CA
INDEX NAME)



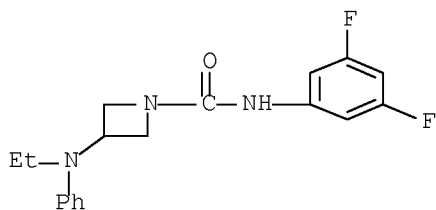
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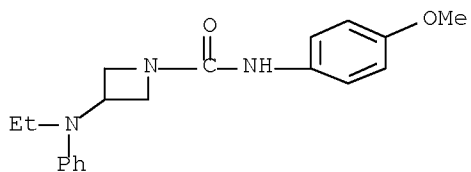
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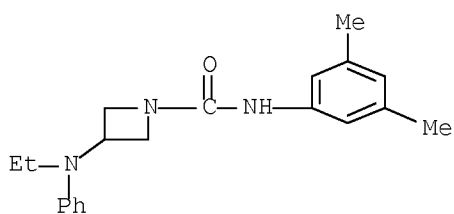
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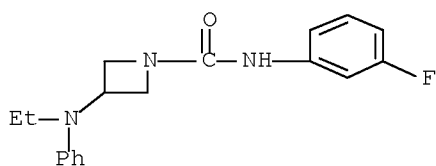
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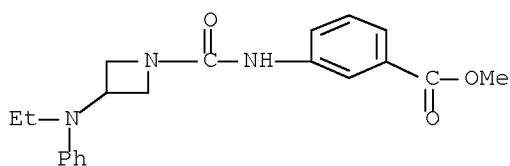
RN 858651-81-7 CAPLUS

CN 1-Azetidinecarboxamide, 3-(ethylphenylamino)-N-(3-fluorophenyl)- (CA INDEX NAME)



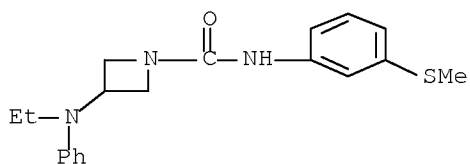
RN 858651-82-8 CAPLUS

CN Benzoic acid, 3-[[[3-(ethylphenylamino)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



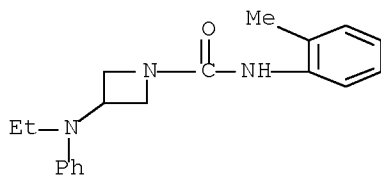
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CN 1-Azetidinecarboxamide, 3-(ethylphenylamino)-N-[3-(methylthio)phenyl]- (CA INDEX NAME)



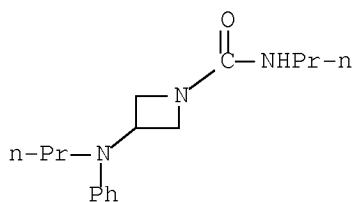
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CN 1-Azetidinecarboxamide, 3-(ethylphenylamino)-N-(2-methylphenyl)- (CA INDEX NAME)



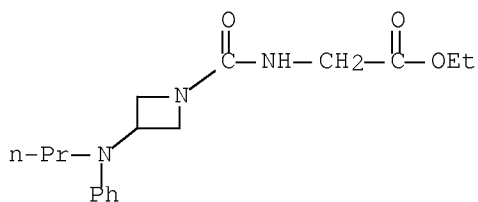
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CN 1-Azetidinecarboxamide, 3-(phenylpropylamino)-N-propyl- (CA INDEX NAME)



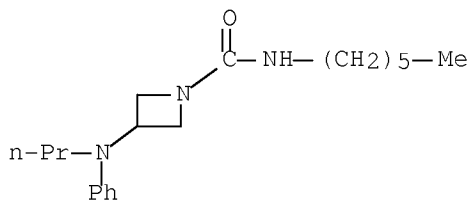
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CN Glycine, N-[[3-(phenylpropylamino)-1-azetidiny]carbonyl]-, ethyl ester
(CA INDEX NAME)



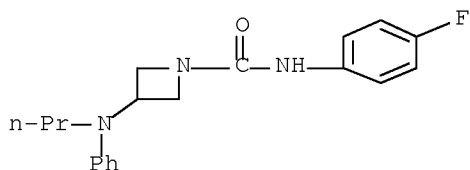
RN 858651-87-3 CAPLUS

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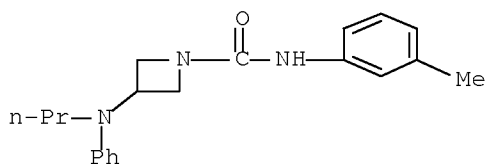
RN 858651-88-4 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-fluorophenyl)-3-(phenylpropylamino)- (CA INDEX NAME)



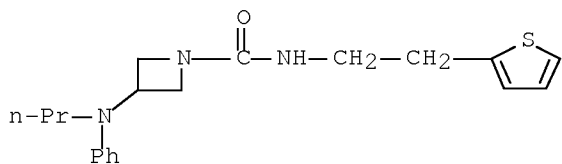
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CN 1-Azetidinecarboxamide, N-(3-methylphenyl)-3-(phenylpropylamino)- (CA INDEX NAME)



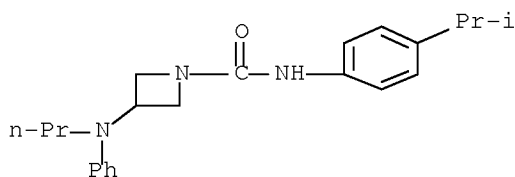
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CN 1-Azetidinecarboxamide, 3-(phenylpropylamino)-N-[2-(2-thienyl)ethyl]- (CA INDEX NAME)



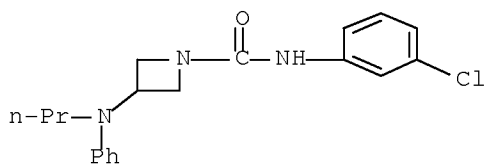
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CN 1-Azetidinecarboxamide, N-[4-(1-methylethyl)phenyl]-3-(phenylpropylamino)- (CA INDEX NAME)



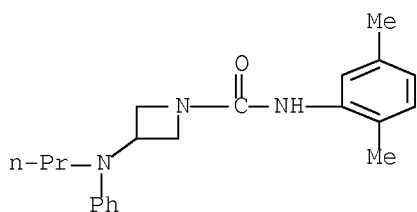
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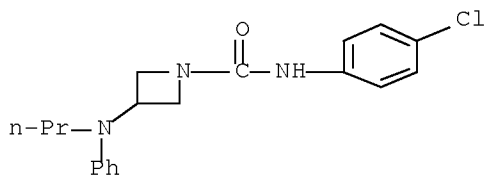
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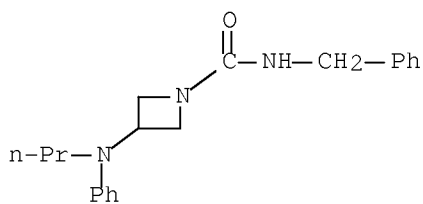
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CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-(phenylpropylamino)- (CA INDEX NAME)



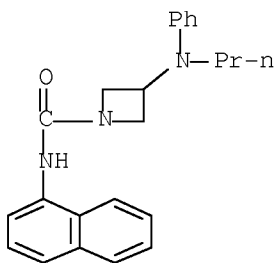
RN 858651-95-3 CAPLUS

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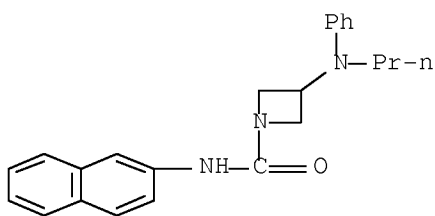
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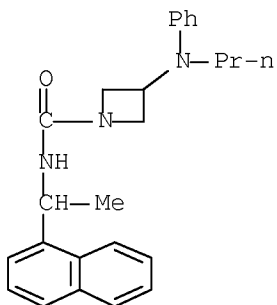
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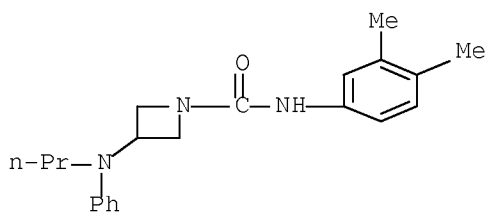
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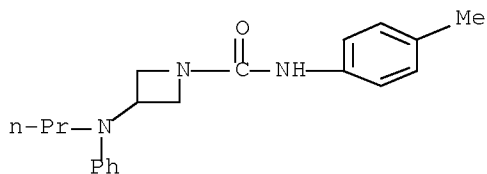
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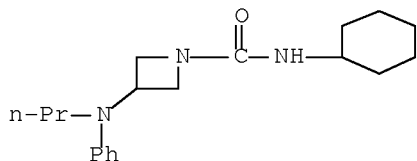
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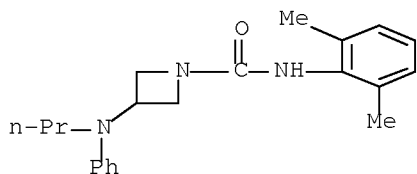
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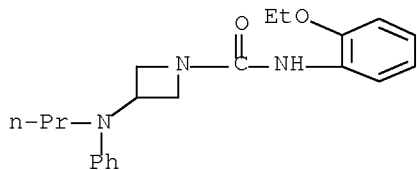
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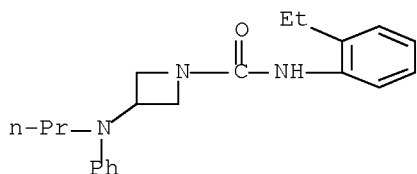


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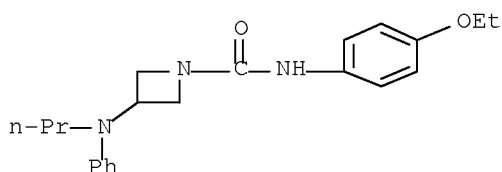
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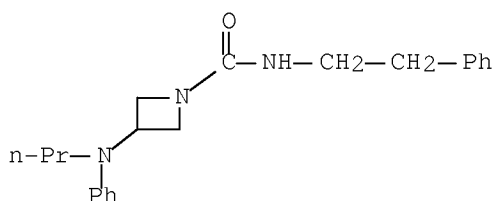
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 INDEX NAME)



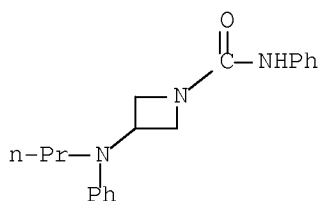
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 CN 1-Azetidinecarboxamide, N-(4-ethoxyphenyl)-3-(phenylpropylamino)- (CA
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 INDEX NAME)

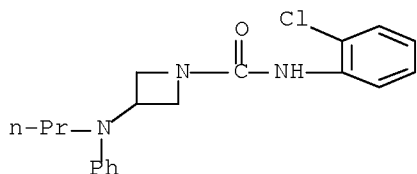


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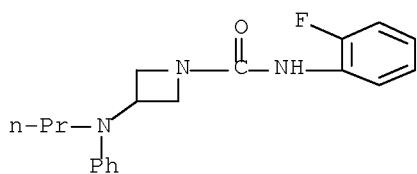
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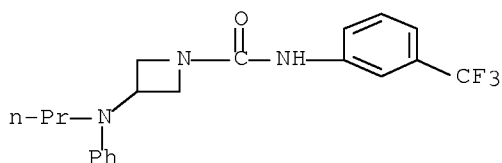
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INDEX NAME)



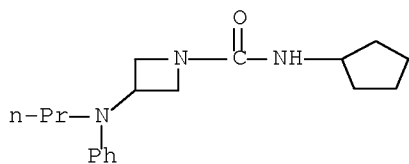
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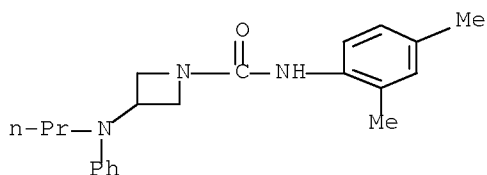
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CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(phenylpropylamino)- (CA INDEX
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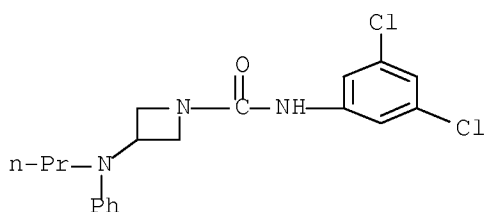
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CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(phenylpropylamino)- (CA
INDEX NAME)



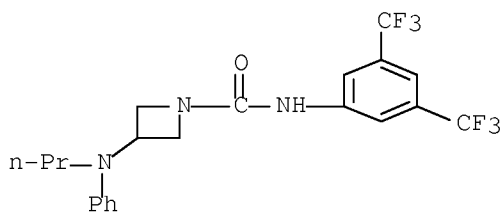
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CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(phenylpropylamino)- (CA INDEX NAME)



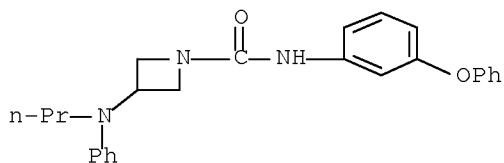
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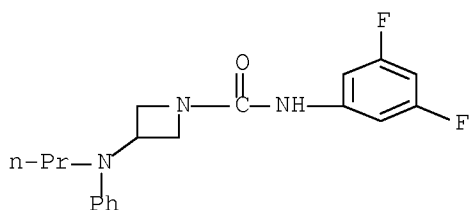
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CN 1-Azetidinecarboxamide, N-(3-phenoxyphenyl)-3-(phenylpropylamino)- (CA INDEX NAME)



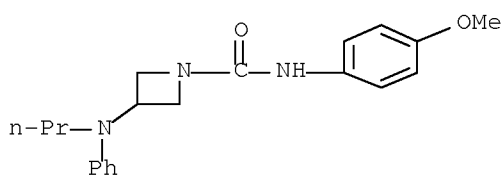
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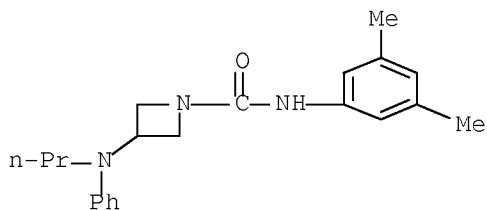
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CN 1-Azetidinecarboxamide, N-(4-methoxyphenyl)-3-(phenylpropylamino)- (CA INDEX NAME)



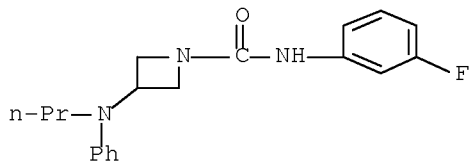
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CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-(phenylpropylamino)- (CA INDEX NAME)



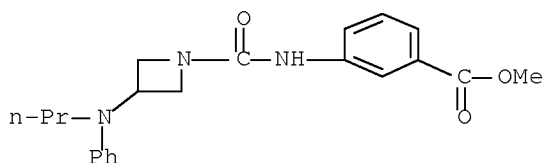
RN 858652-19-4 CAPLUS

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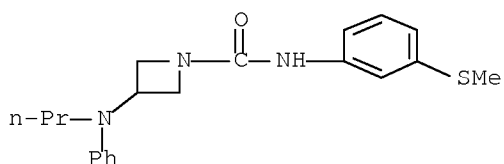
RN 858652-20-7 CAPLUS

CN Benzoic acid, 3-[[[3-(phenylpropylamino)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



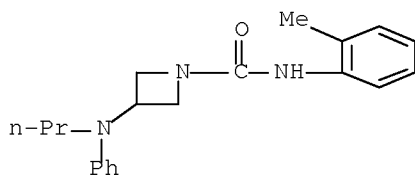
RN 858652-21-8 CAPLUS

CN 1-Azetidinecarboxamide, N-[3-(methylthio)phenyl]-3-(phenylpropylamino)-
(CA INDEX NAME)



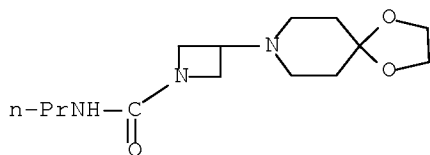
RN 858652-22-9 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-methylphenyl)-3-(phenylpropylamino)- (CA
INDEX NAME)



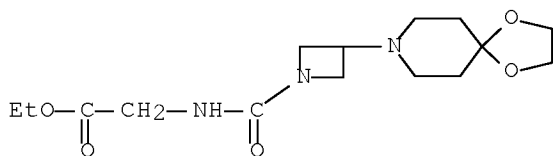
RN 858652-23-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-propyl-
(CA INDEX NAME)



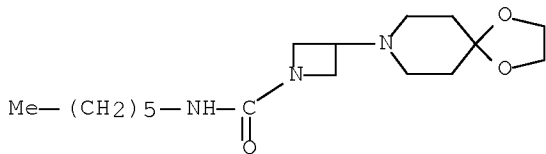
RN 858652-24-1 CAPLUS

CN Glycine, N-[[3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-1-azetidiny]carbonyl]-
, ethyl ester (CA INDEX NAME)



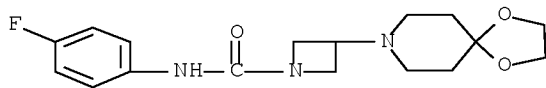
RN 858652-25-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-hexyl-
(CA INDEX NAME)



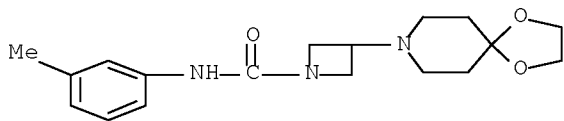
RN 858652-26-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-(4-fluorophenyl)-
(CA INDEX NAME)



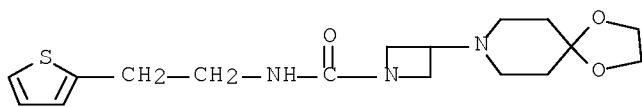
RN 858652-27-4 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-(3-methylphenyl)-
(CA INDEX NAME)



RN 858652-28-5 CAPLUS

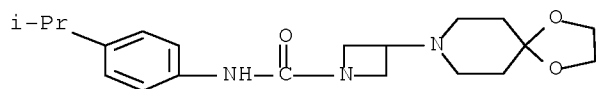
CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-[2-(2-thienyl)ethyl]-
(CA INDEX NAME)



RN 858652-29-6 CAPLUS

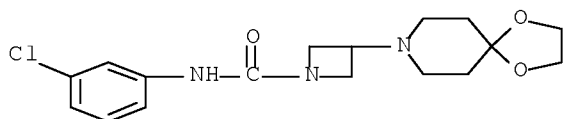
CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-[4-(1-methyl-2-thienyl)ethyl]-
(CA INDEX NAME)

methylethyl)phenyl]- (CA INDEX NAME)



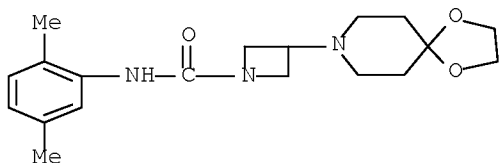
RN 858652-30-9 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-(1,4-dioxaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



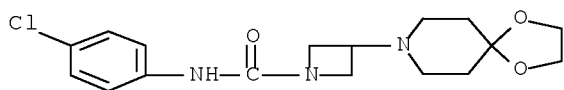
RN 858652-31-0 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,5-dimethylphenyl)-3-(1,4-dioxaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



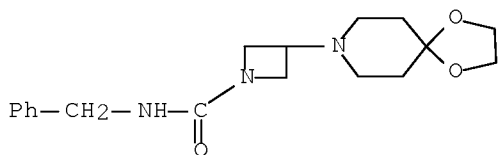
RN 858652-32-1 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-(1,4-dioxaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



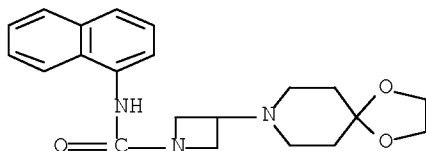
RN 858652-34-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxaspiro[4.5]dec-8-yl)-N-(phenylmethyl)- (CA INDEX NAME)



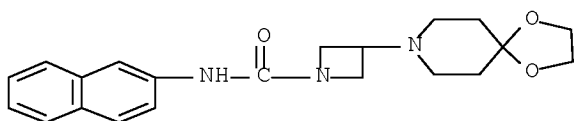
RN 858652-36-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-1-naphthalenyl- (CA INDEX NAME)



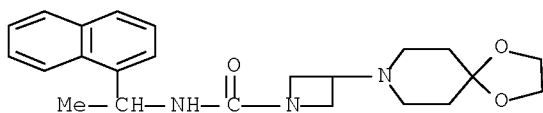
RN 858652-37-6 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-2-naphthalenyl- (CA INDEX NAME)



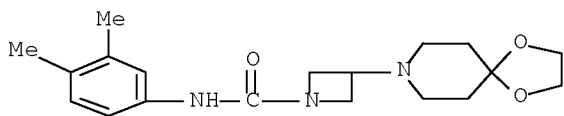
RN 858652-38-7 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-[1-(1-naphthalenyl)ethyl]- (CA INDEX NAME)



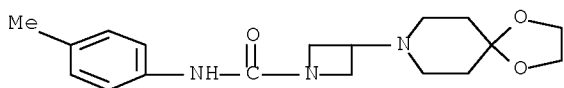
RN 858652-39-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,4-dimethylphenyl)-3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



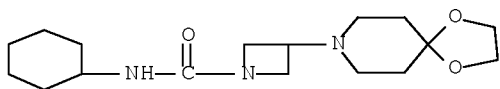
RN 858652-40-1 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-(4-methylphenyl)- (CA INDEX NAME)



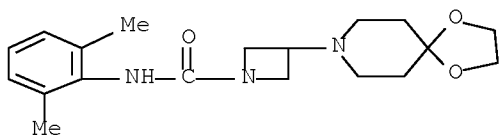
RN 858652-41-2 CAPLUS

CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(1,4-dioxaspiro[4.5]dec-8-yl)-
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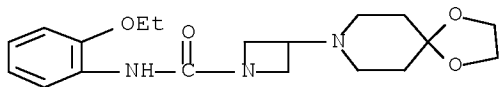
RN 858652-42-3 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-(1,4-dioxaspiro[4.5]dec-8-yl)-
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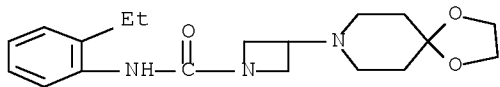
RN 858652-43-4 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxaspiro[4.5]dec-8-yl)-N-(2-ethoxyphenyl)-
(CA INDEX NAME)



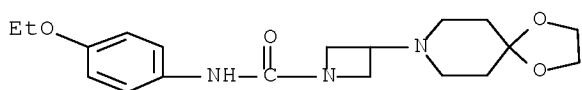
RN 858652-44-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxaspiro[4.5]dec-8-yl)-N-(2-ethylphenyl)-
(CA INDEX NAME)



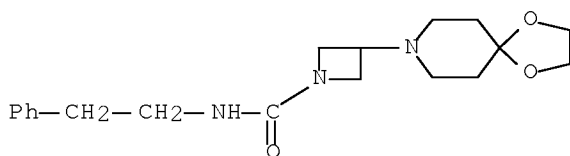
RN 858652-45-6 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxaspiro[4.5]dec-8-yl)-N-(4-ethoxyphenyl)-
(CA INDEX NAME)



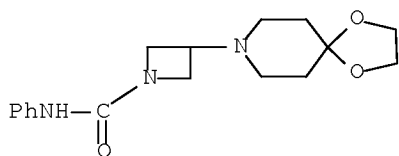
RN 858652-46-7 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-(2-phenylethyl)- (CA INDEX NAME)



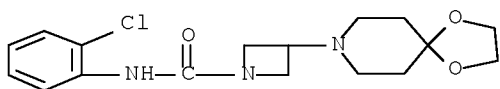
RN 858652-47-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-phenyl- (CA INDEX NAME)



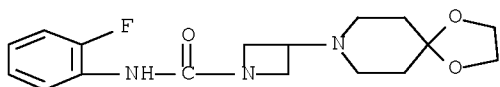
RN 858652-48-9 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-chlorophenyl)-3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



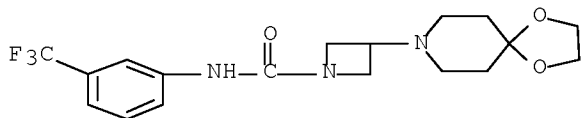
RN 858652-49-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-(2-fluorophenyl)- (CA INDEX NAME)



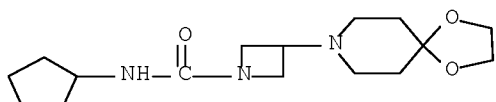
RN 858652-50-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



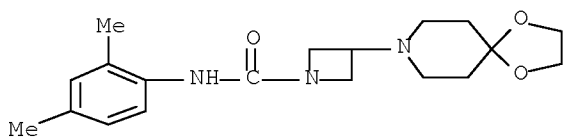
RN 858652-51-4 CAPLUS

CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(1,4-dioxaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



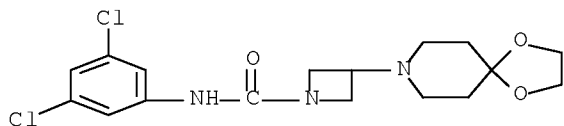
RN 858652-52-5 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(1,4-dioxaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



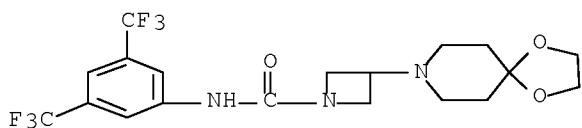
RN 858652-53-6 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(1,4-dioxaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



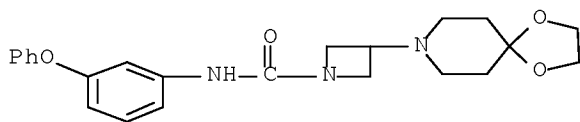
RN 858652-54-7 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(1,4-dioxaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



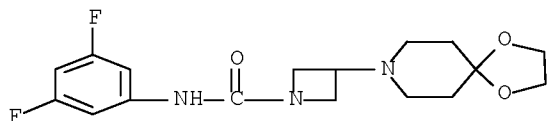
RN 858652-55-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-(3-phenoxyphenyl)- (CA INDEX NAME)



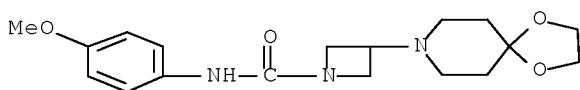
RN 858652-56-9 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



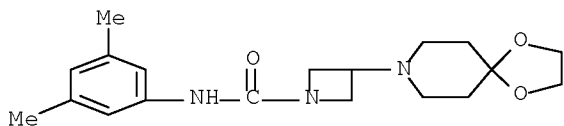
RN 858652-57-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-(4-methoxyphenyl)- (CA INDEX NAME)



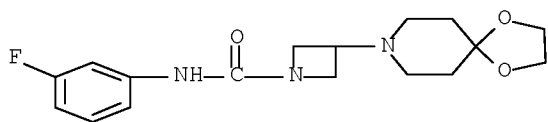
RN 858652-58-1 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)- (CA INDEX NAME)



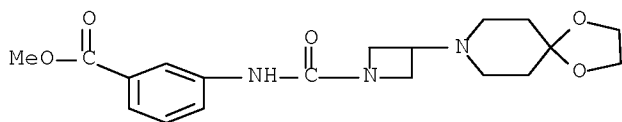
RN 858652-59-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-(3-fluorophenyl)- (CA INDEX NAME)



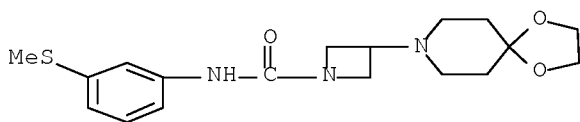
RN 858652-60-5 CAPLUS

CN Benzoic acid, 3-[[[3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



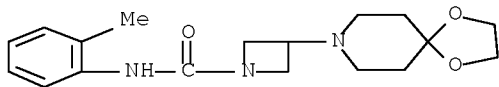
RN 858652-61-6 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-[3-(methylthio)phenyl]- (CA INDEX NAME)



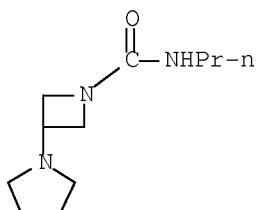
RN 858652-62-7 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1,4-dioxo-8-azaspiro[4.5]dec-8-yl)-N-(2-methylphenyl)- (CA INDEX NAME)



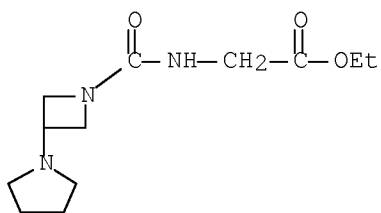
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CN 1-Azetidinecarboxamide, N-propyl-3-(1-pyrrolidinyl)- (CA INDEX NAME)



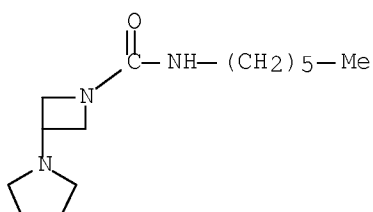
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CN Glycine, N-[[3-(1-pyrrolidinyl)-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)



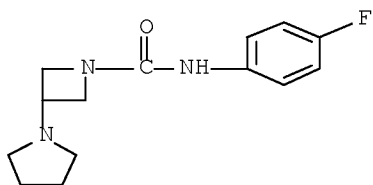
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CN 1-Azetidinecarboxamide, N-hexyl-3-(1-pyrrolidinyl)- (CA INDEX NAME)



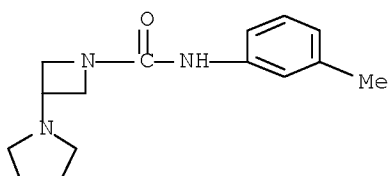
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CN 1-Azetidinecarboxamide, N-(4-fluorophenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



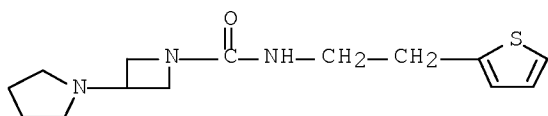
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CN 1-Azetidinecarboxamide, N-(3-methylphenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



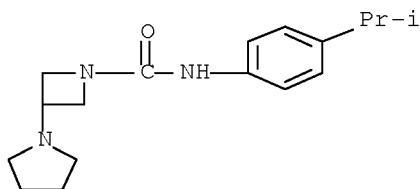
RN 858652-68-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(1-pyrrolidinyl)-N-[2-(2-thienyl)ethyl]- (CA INDEX NAME)



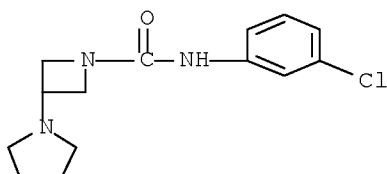
RN 858652-69-4 CAPLUS

CN 1-Azetidinecarboxamide, N-[4-(1-methylethyl)phenyl]-3-(1-pyrrolidinyl)-
(CA INDEX NAME)



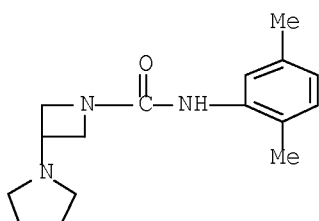
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CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



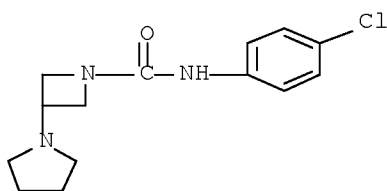
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CN 1-Azetidinecarboxamide, N-(2,5-dimethylphenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)

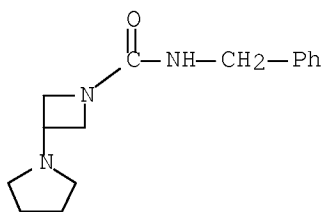


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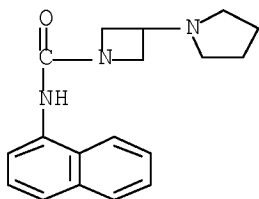
CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



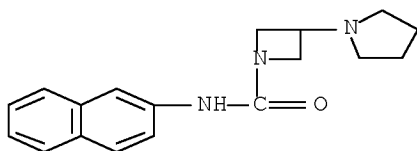
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 CN 1-Azetidinecarboxamide, N-(phenylmethyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



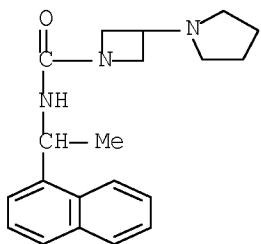
RN 858652-74-1 CAPLUS
 CN 1-Azetidinecarboxamide, N-1-naphthalenyl-3-(1-pyrrolidinyl)- (CA INDEX NAME)



RN 858652-75-2 CAPLUS
 CN 1-Azetidinecarboxamide, N-2-naphthalenyl-3-(1-pyrrolidinyl)- (CA INDEX NAME)

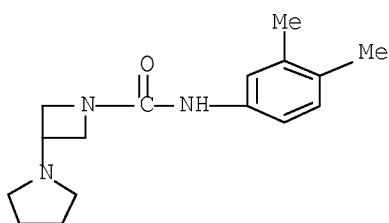


RN 858652-76-3 CAPLUS
 CN 1-Azetidinecarboxamide, N-[1-(1-naphthalenyl)ethyl]-3-(1-pyrrolidinyl)- (CA INDEX NAME)



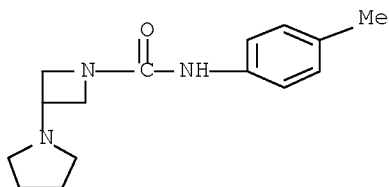
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CN 1-Azetidinecarboxamide, N-(3,4-dimethylphenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



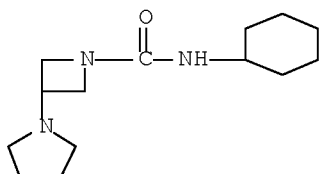
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CN 1-Azetidinecarboxamide, N-(4-methylphenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



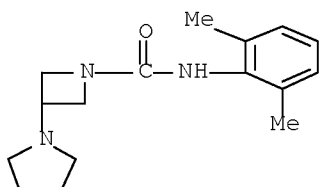
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CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(1-pyrrolidinyl)- (CA INDEX NAME)



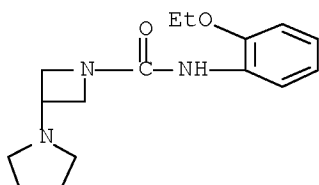
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CN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



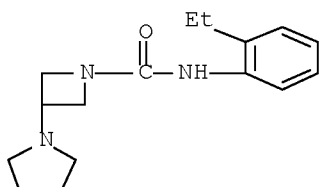
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CN 1-Azetidinecarboxamide, N-(2-ethoxyphenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



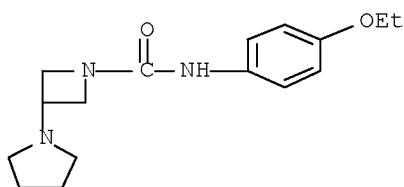
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CN 1-Azetidinecarboxamide, N-(2-ethylphenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



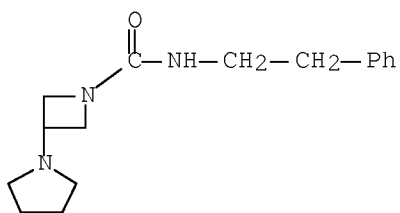
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CN 1-Azetidinecarboxamide, N-(4-ethoxyphenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



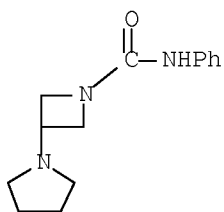
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CN 1-Azetidinecarboxamide, N-(2-phenylethyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



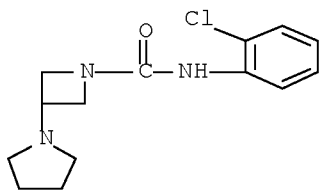
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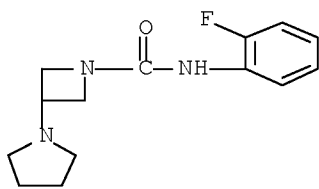
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CN 1-Azetidinecarboxamide, N-(2-chlorophenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



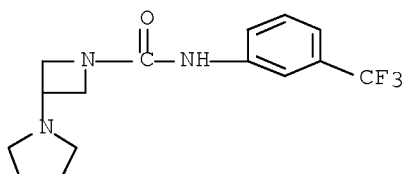
RN 858652-87-6 CAPLUS

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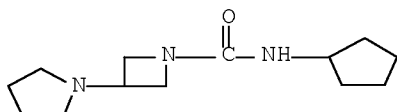
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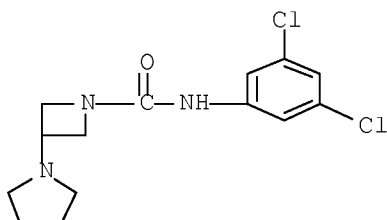
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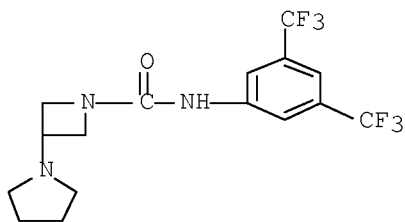
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CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



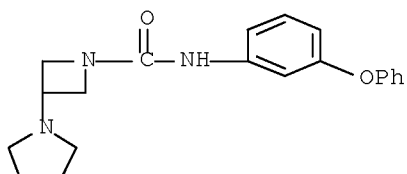
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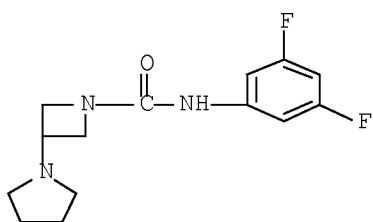
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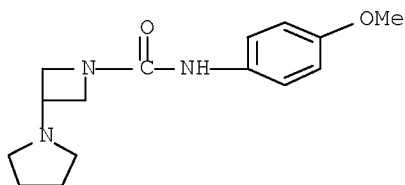
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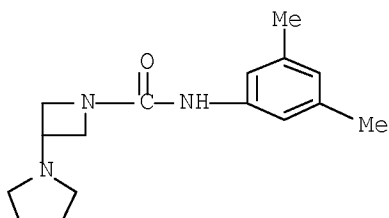
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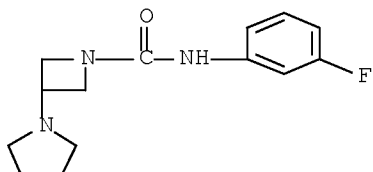
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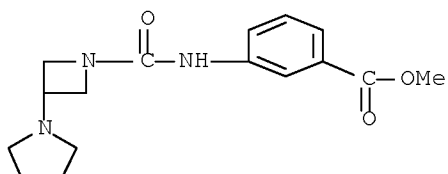
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CN 1-Azetidinecarboxamide, N-(3-fluorophenyl)-3-(1-pyrrolidinyl)- (CA INDEX NAME)



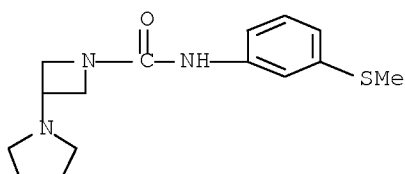
RN 858652-97-8 CAPLUS

CN Benzoic acid, 3-[[[3-(1-pyrrolidinyl)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



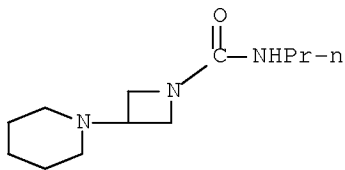
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CN 1-Azetidinecarboxamide, N-[3-(methylthio)phenyl]-3-(1-pyrrolidinyl)- (CA INDEX NAME)



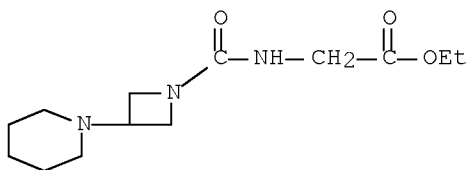
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CN 1-Azetidinecarboxamide, 3-(1-piperidinyl)-N-propyl- (CA INDEX NAME)



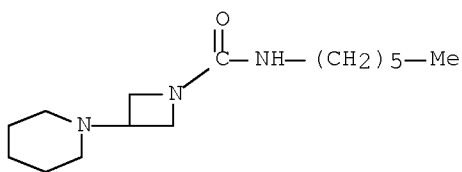
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CN Glycine, N-[[[3-(1-piperidinyl)-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)



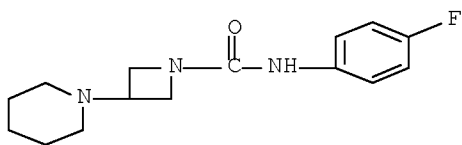
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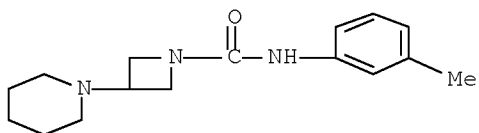
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CN 1-Azetidinecarboxamide, N-(4-fluorophenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



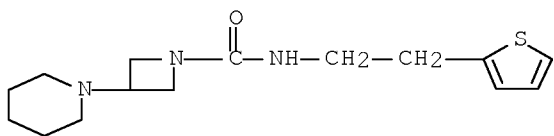
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CN 1-Azetidinecarboxamide, N-(3-methylphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



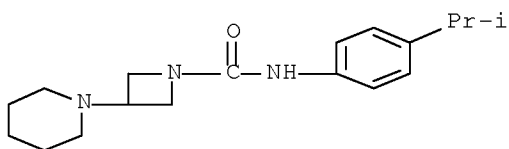
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CN 1-Azetidinecarboxamide, 3-(1-piperidinyl)-N-[2-(2-thienyl)ethyl]- (CA INDEX NAME)



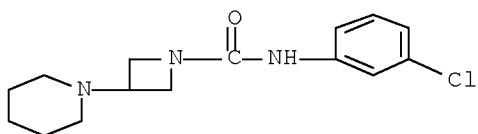
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CN 1-Azetidinecarboxamide, N-[4-(1-methylethyl)phenyl]-3-(1-piperidinyl)-
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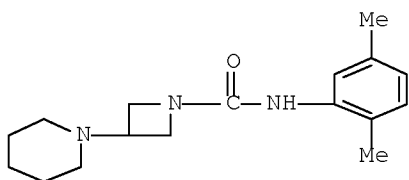
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CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-(1-piperidinyl)- (CA INDEX
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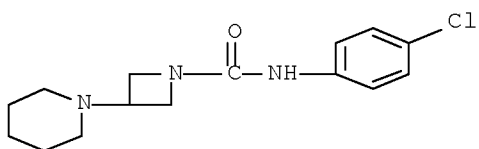
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CN 1-Azetidinecarboxamide, N-(2,5-dimethylphenyl)-3-(1-piperidinyl)- (CA
INDEX NAME)



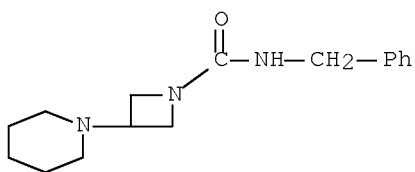
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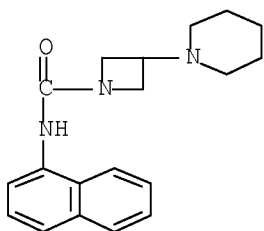
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CN 1-Azetidinecarboxamide, N-(phenylmethyl)-3-(1-piperidinyl)- (CA INDEX NAME)



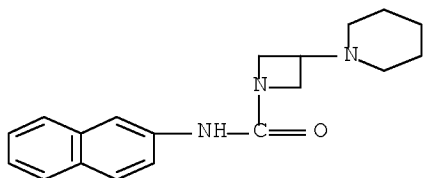
RN 858653-10-8 CAPLUS

CN 1-Azetidinecarboxamide, N-1-naphthalenyl-3-(1-piperidinyl)- (CA INDEX NAME)



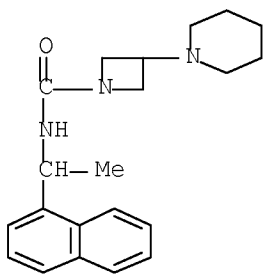
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CN 1-Azetidinecarboxamide, N-2-naphthalenyl-3-(1-piperidinyl)- (CA INDEX NAME)



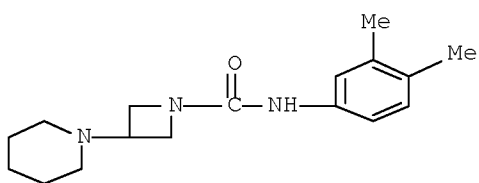
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CN 1-Azetidinecarboxamide, N-[1-(1-naphthalenyl)ethyl]-3-(1-piperidinyl)- (CA INDEX NAME)



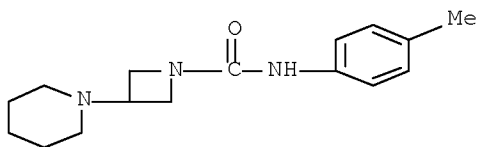
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CN 1-Azetidinecarboxamide, N-(3,4-dimethylphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



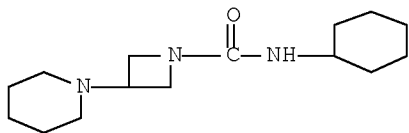
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CN 1-Azetidinecarboxamide, N-(4-methylphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



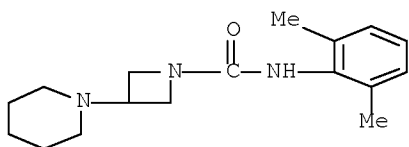
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CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(1-piperidinyl)- (CA INDEX NAME)



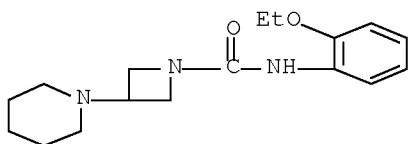
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CN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



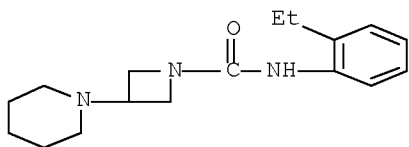
RN 858653-17-5 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-ethoxyphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



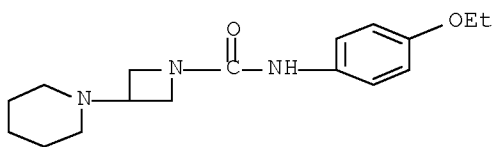
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CN 1-Azetidinecarboxamide, N-(2-ethylphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



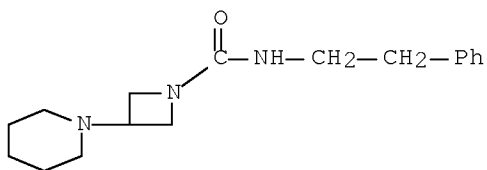
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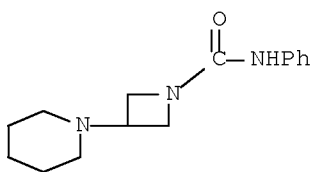
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CN 1-Azetidinecarboxamide, N-(2-phenylethyl)-3-(1-piperidinyl)- (CA INDEX NAME)



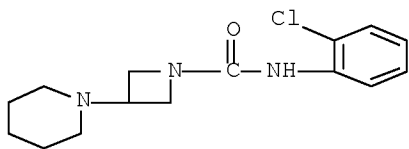
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CN 1-Azetidinecarboxamide, N-phenyl-3-(1-piperidinyl)- (CA INDEX NAME)



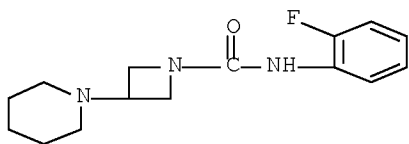
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CN 1-Azetidinecarboxamide, N-(2-chlorophenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



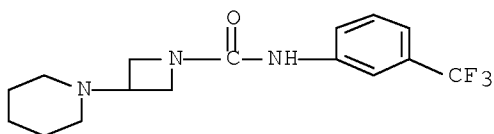
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CN 1-Azetidinecarboxamide, N-(2-fluorophenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



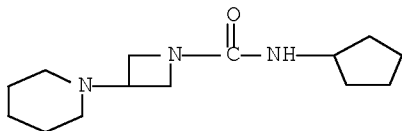
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CN 1-Azetidinecarboxamide, 3-(1-piperidinyl)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



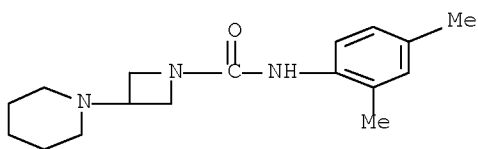
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CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(1-piperidinyl)- (CA INDEX NAME)



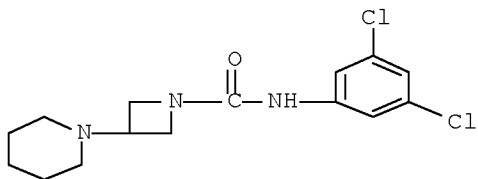
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CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



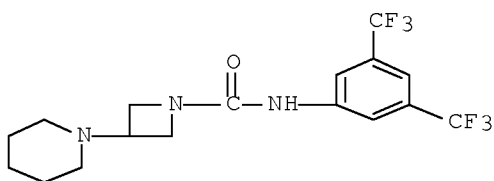
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CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(1-piperidinyl)- (CA INDEX NAME)

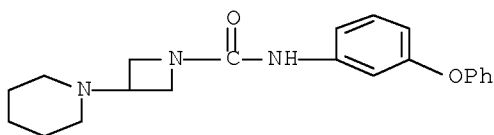


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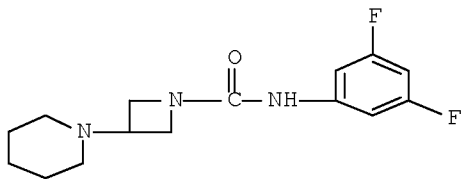
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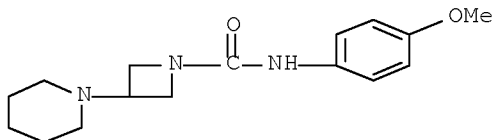
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 CN 1-Azetidinecarboxamide, N-(3-phenoxyphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



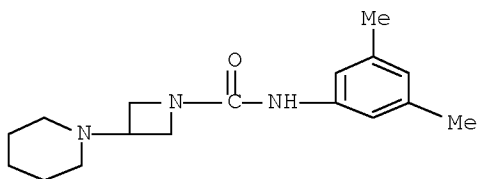
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 CN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



RN 858653-31-3 CAPLUS
 CN 1-Azetidinecarboxamide, N-(4-methoxyphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)

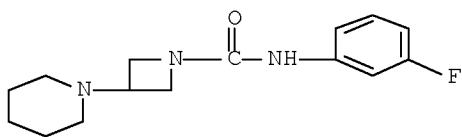


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 CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



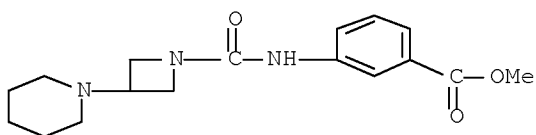
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CN 1-Azetidinecarboxamide, N-(3-fluorophenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



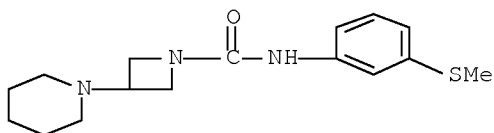
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CN Benzoic acid, 3-[[[3-(1-piperidinyl)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



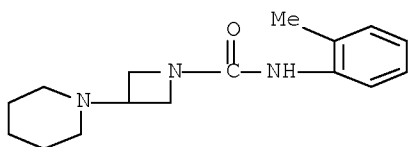
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CN 1-Azetidinecarboxamide, N-[3-(methylthio)phenyl]-3-(1-piperidinyl)- (CA INDEX NAME)



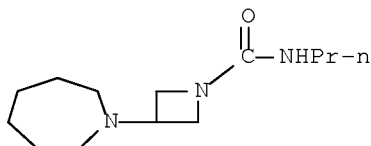
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CN 1-Azetidinecarboxamide, N-(2-methylphenyl)-3-(1-piperidinyl)- (CA INDEX NAME)



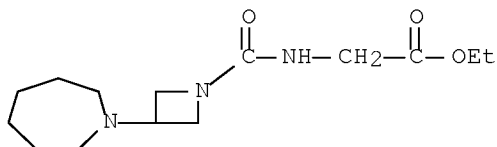
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CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-propyl- (CA INDEX NAME)



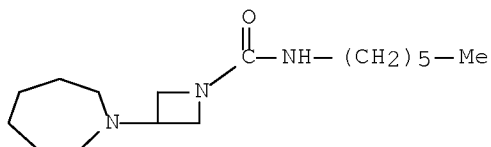
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CN Glycine, N-[[3-(hexahydro-1H-azepin-1-yl)-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)



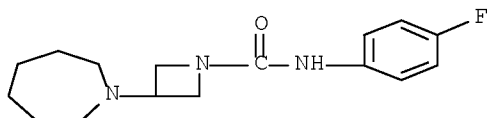
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CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-hexyl- (CA INDEX NAME)

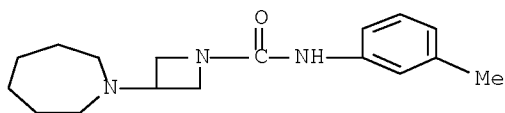


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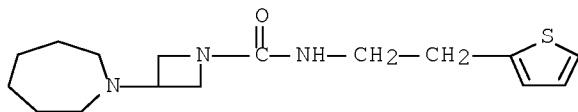
CN 1-Azetidinecarboxamide, N-(4-fluorophenyl)-3-(hexahydro-1H-azepin-1-yl)- (CA INDEX NAME)



RN 858653-41-5 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-(3-methylphenyl)-
 (CA INDEX NAME)



RN 858653-42-6 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-[2-(2-thienyl)ethyl]- (CA INDEX NAME)



IT 858653-43-7P, 3-(Azepan-1-yl)-N-(4-isopropylphenyl)azetidine-1-carboxamide 858653-44-8P, 3-(Azepan-1-yl)-N-(3-chlorophenyl)azetidine-1-carboxamide 858653-45-9P, 3-(Azepan-1-yl)-N-(2,5-dimethylphenyl)azetidine-1-carboxamide 858653-46-0P, 3-(Azepan-1-yl)-N-(4-chlorophenyl)azetidine-1-carboxamide 858653-47-1P, 3-(Azepan-1-yl)-N-benzylazetidine-1-carboxamide 858653-48-2P, 3-(Azepan-1-yl)-N-(1-naphthyl)azetidine-1-carboxamide 858653-49-3P, 3-(Azepan-1-yl)-N-(2-naphthyl)azetidine-1-carboxamide 858653-50-6P, 3-(Azepan-1-yl)-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858653-51-7P, 3-(Azepan-1-yl)-N-(3,4-dimethylphenyl)azetidine-1-carboxamide 858653-52-8P, 3-(Azepan-1-yl)-N-(4-methylphenyl)azetidine-1-carboxamide 858653-53-9P, 3-(Azepan-1-yl)-N-cyclohexylazetidine-1-carboxamide 858653-54-0P, 3-(Azepan-1-yl)-N-(2,6-dimethylphenyl)azetidine-1-carboxamide 858653-55-1P, 3-(Azepan-1-yl)-N-(2-ethoxyphenyl)azetidine-1-carboxamide 858653-56-2P, 3-(Azepan-1-yl)-N-(2-ethylphenyl)azetidine-1-carboxamide 858653-57-3P, 3-(Azepan-1-yl)-N-(4-ethoxyphenyl)azetidine-1-carboxamide 858653-58-4P, 3-(Azepan-1-yl)-N-(2-phenylethyl)azetidine-1-carboxamide 858653-59-5P, 3-(Azepan-1-yl)-N-phenylazetidine-1-carboxamide 858653-60-8P, 3-(Azepan-1-yl)-N-(2-chlorophenyl)azetidine-1-carboxamide 858653-61-9P, 3-(Azepan-1-yl)-N-(2-fluorophenyl)azetidine-1-carboxamide 858653-62-0P, 3-(Azepan-1-yl)-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858653-63-1P, 3-(Azepan-1-yl)-N-cyclopentylazetidine-1-carboxamide 858653-64-2P, 3-(Azepan-1-yl)-N-(2,4-dimethylphenyl)azetidine-1-carboxamide 858653-65-3P, 3-(Azepan-1-yl)-N-(3,5-dichlorophenyl)azetidine-1-carboxamide 858653-66-4P, 3-(Azepan-1-yl)-N-[3,5-bis(trifluoromethyl)phenyl]azetidine-1-carboxamide 858653-67-5P, 3-(Azepan-1-yl)-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858653-68-6P, 3-(Azepan-1-yl)-N-(3,5-difluorophenyl)azetidine-1-carboxamide 858653-69-7P, 3-(Azepan-1-yl)-N-(4-

methoxyphenyl)azetidine-1-carboxamide 858653-70-0P,
3-(Azepan-1-yl)-N-(3,5-dimethylphenyl)azetidine-1-carboxamide
858653-71-1P, 3-(Azepan-1-yl)-N-(3-fluorophenyl)azetidine-1-
carboxamide 858653-72-2P, Methyl 3-[[3-(azepan-1-yl)azetidin-1-
yl]carbonyl]amino]benzoate 858653-73-3P, 3-(Azepan-1-yl)-N-[3-
(methylsulfanyl)phenyl]azetidine-1-carboxamide 858653-74-4P,
3-(Azepan-1-yl)-N-(2-methylphenyl)azetidine-1-carboxamide
858653-75-5P, 3-(Diisobutylamino)-N-ethylazetidine-1-carboxamide
858653-76-6P, 3-(Diisobutylamino)-N-propylazetidine-1-carboxamide
858653-77-7P 858653-78-8P, 3-(Diisobutylamino)-N-
hexylazetidine-1-carboxamide 858653-79-9P, 3-(Diisobutylamino)-N-
(4-fluorophenyl)azetidine-1-carboxamide 858653-80-2P,
3-(Diisobutylamino)-N-(3-methylphenyl)azetidine-1-carboxamide
858653-81-3P, 3-(Diisobutylamino)-N-(2-(thien-2-yl)ethyl)azetidine-
1-carboxamide 858653-82-4P, 3-(Diisobutylamino)-N-(4-
isopropylphenyl)azetidine-1-carboxamide 858653-83-5P,
N-(3-Chlorophenyl)-3-(diisobutylamino)azetidine-1-carboxamide
858653-84-6P, 3-(Diisobutylamino)-N-(2,5-dimethylphenyl)azetidine-
1-carboxamide 858653-85-7P, N-(4-Chlorophenyl)-3-
(diisobutylamino)azetidine-1-carboxamide 858653-86-8P,
N-Benzyl-3-(diisobutylamino)azetidine-1-carboxamide 858653-87-9P
, 3-(Diisobutylamino)-N-(1-naphthyl)azetidine-1-carboxamide
858653-88-0P, 3-(Diisobutylamino)-N-(2-naphthyl)azetidine-1-
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naphthyl)ethyl]azetidine-1-carboxamide 858653-90-4P,
3-(Diisobutylamino)-N-(3,4-dimethylphenyl)azetidine-1-carboxamide
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carboxamide 858653-92-6P, N-Cyclohexyl-3-
(diisobutylamino)azetidine-1-carboxamide 858653-93-7P,
3-(Diisobutylamino)-N-(2,6-dimethylphenyl)azetidine-1-carboxamide
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3-(Diisobutylamino)-N-(4-ethoxyphenyl)azetidine-1-carboxamide
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carboxamide 858653-98-2P, 3-(Diisobutylamino)-N-phenylazetidine-
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(diisobutylamino)azetidine-1-carboxamide 858654-00-9P,
3-(Diisobutylamino)-N-(2-fluorophenyl)azetidine-1-carboxamide
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(trifluoromethyl)phenyl]azetidine-1-carboxamide 858654-02-1P,
N-Cyclopentyl-3-(diisobutylamino)azetidine-1-carboxamide
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(diisobutylamino)azetidine-1-carboxamide 858654-05-4P,
N-[3,5-Bis(trifluoromethyl)phenyl]-3-(diisobutylamino)azetidine-1-
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phenoxyphenyl)azetidine-1-carboxamide 858654-07-6P,
N-(3,5-Difluorophenyl)-3-(diisobutylamino)azetidine-1-carboxamide
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carboxamide 858654-09-8P, 3-(Diisobutylamino)-N-(3,5-
dimethylphenyl)azetidine-1-carboxamide 858654-10-1P,
3-(Diisobutylamino)-N-(3-fluorophenyl)azetidine-1-carboxamide
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yl]carbonyl]amino)benzoate 858654-12-3P, 3-(Diisobutylamino)-N-
[3-(methylsulfanyl)phenyl]azetidine-1-carboxamide 858654-13-4P,
3-(Diisobutylamino)-N-(2-methylphenyl)azetidine-1-carboxamide
858654-14-5P, 3-Morpholin-4-yl-N-propylazetidine-1-carboxamide
858654-15-6P, N-Hexyl-3-morpholin-4-ylazetidine-1-carboxamide
858654-16-7P, N-(4-Fluorophenyl)-3-morpholin-4-ylazetidine-1-

carboxamide 858654-17-8P, N-(3-Methylphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-18-9P, 3-Morpholin-4-yl-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858654-19-0P, N-(4-Isopropylphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-20-3P, N-(3-Chlorophenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-21-4P, N-(2,5-Dimethylphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-22-5P, N-(4-Chlorophenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-23-6P, N-Benzyl-3-morpholin-4-ylazetidine-1-carboxamide 858654-24-7P, 3-Morpholin-4-yl-N-(1-naphthyl)azetidine-1-carboxamide 858654-25-8P, 3-Morpholin-4-yl-N-(2-naphthyl)azetidine-1-carboxamide 858654-27-0P, 3-Morpholin-4-yl-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858654-28-1P, N-(3,4-Dimethylphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-30-5P, N-(4-Methylphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-31-6P, N-Cyclohexyl-3-morpholin-4-ylazetidine-1-carboxamide 858654-32-7P, N-(2,6-Dimethylphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-33-8P, N-(2-Ethoxyphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-34-9P, N-(2-Ethylphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-35-0P, N-(4-Ethoxyphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-36-1P, 3-Morpholin-4-yl-N-(2-phenylethyl)azetidine-1-carboxamide 858654-37-2P, 3-Morpholin-4-yl-N-phenylazetidine-1-carboxamide 858654-38-3P, N-(2-Chlorophenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-39-4P, N-(2-Fluorophenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-40-7P, 3-Morpholin-4-yl-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858654-41-8P, N-Cyclopentyl-3-morpholin-4-ylazetidine-1-carboxamide 858654-42-9P, N-(2,4-Dimethylphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-43-0P, N-(3,5-Dichlorophenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-44-1P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-morpholin-4-ylazetidine-1-carboxamide 858654-45-2P, 3-Morpholin-4-yl-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858654-46-3P, N-(3,5-Difluorophenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-47-4P, N-(4-Methoxyphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-48-5P, N-(3,5-Dimethylphenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-49-6P, N-(3-Fluorophenyl)-3-morpholin-4-ylazetidine-1-carboxamide 858654-50-9P, Methyl 3-[[3-morpholin-4-ylazetidin-1-yl)carbonyl]amino]benzoate 858654-51-0P, N-[3-(Methylsulfanyl)phenyl]-3-morpholin-4-ylazetidine-1-carboxamide 858654-52-1P, N-Hexyl-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-53-2P, N-(4-Fluorophenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-54-3P, N-(3-Methylphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-55-4P, 3-(4-Methylpiperazin-1-yl)-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858654-56-5P, N-(4-Isopropylphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-57-6P, N-(3-Chlorophenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-58-7P, N-(2,5-Dimethylphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-59-8P, N-(4-Chlorophenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-60-1P, N-Benzyl-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-61-2P, 3-(4-Methylpiperazin-1-yl)-N-(1-naphthyl)azetidine-1-carboxamide 858654-62-3P, 3-(4-Methylpiperazin-1-yl)-N-(2-naphthyl)azetidine-1-carboxamide 858654-63-4P, 3-(4-Methylpiperazin-1-yl)-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858654-64-5P, N-(3,4-Dimethylphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide

858654-65-6P, N-(4-Methylphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-66-7P, N-Cyclohexyl-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-67-8P, N-(2,6-Dimethylphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-68-9P, N-(2-Ethoxyphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-69-0P, N-(4-Ethoxyphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-70-3P, 3-(4-Methylpiperazin-1-yl)-N-(2-phenylethyl)azetidine-1-carboxamide 858654-71-4P, 3-(4-Methylpiperazin-1-yl)-N-phenylazetidine-1-carboxamide 858654-72-5P, N-(2-Chlorophenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-73-6P, N-(2-Fluorophenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-74-7P, N-Cyclopentyl-3-(4-Methylpiperazin-1-yl)azetidine-1-carboxamide 858654-75-8P, N-(3,5-Dichlorophenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-76-9P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-77-0P, 3-(4-Methylpiperazin-1-yl)-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858654-78-1P, N-(3,5-Difluorophenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-79-2P, N-(4-Methoxyphenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-80-5P, N-(3-Fluorophenyl)-3-(4-methylpiperazin-1-yl)azetidine-1-carboxamide 858654-81-6P, 3-(4-Methylpiperazin-1-yl)-N-[3-(methylsulfonyl)phenyl]azetidine-1-carboxamide 858654-82-7P, N-Ethyl-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-83-8P, 3-(4-Phenylpiperazin-1-yl)-N-propylazetidine-1-carboxamide 858654-84-9P 858654-85-0P, N-Hexyl-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-86-1P, N-(4-Fluorophenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-87-2P, N-(3-Methylphenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-88-3P, 3-(4-Phenylpiperazin-1-yl)-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858654-89-4P, N-(4-Isopropylphenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-90-7P, N-(3-Chlorophenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-91-8P, N-(2,5-Dimethylphenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-92-9P, N-(4-Chlorophenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-93-0P, N-Benzyl-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-94-1P, N-(1-Naphthyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-97-4P, N-(2-Naphthyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-98-5P, N-[1-(1-Naphthyl)ethyl]-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858654-99-6P, N-(3,4-Dimethylphenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-00-2P, N-(4-Methylphenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-01-3P, N-Cyclohexyl-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-02-4P, N-(2,6-Dimethylphenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-03-5P, N-(2-Ethoxyphenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-04-6P, N-(2-Ethylphenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-05-7P, N-(4-Ethoxyphenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-06-8P, N-(2-Phenylethyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-07-9P, N-Phenyl-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-08-0P, N-(2-Chlorophenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-09-1P, N-(2-Fluorophenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-10-4P, 3-(4-Phenylpiperazin-1-yl)-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858655-11-5P, N-Cyclopentyl-3-(4-Phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-12-6P, N-(2,4-Dimethylphenyl)-

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858655-14-8P, N-(3-Phenoxyphenyl)-3-(4-phenylpiperazin-1-
yl)azetidine-1-carboxamide 858655-15-9P, N-(3,5-Difluorophenyl)-
3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-16-0P,
N-(4-Methoxyphenyl)-3-(4-phenylpiperazin-1-yl)azetidine-1-carboxamide
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yl)azetidine-1-carboxamide 858655-18-2P, N-(3-Fluorophenyl)-3-(4-
phenylpiperazin-1-yl)azetidine-1-carboxamide 858655-19-3P,
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yl)azetidine-1-carboxamide 858655-22-8P, N-Ethyl-3-(4-
phenylpiperidin-1-yl)azetidine-1-carboxamide 858655-23-9P,
3-(4-Phenylpiperidin-1-yl)-N-propylazetidine-1-carboxamide
858655-24-0P 858655-25-1P, N-Hexyl-3-(4-phenylpiperidin-
1-yl)azetidine-1-carboxamide 858655-26-2P, N-(4-Fluorophenyl)-3-
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N-(3-Methylphenyl)-3-(4-phenylpiperidin-1-yl)azetidine-1-carboxamide
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yl)ethyl)azetidine-1-carboxamide 858655-29-5P,
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858655-57-9P, Methyl 3-([3-(4-phenylpiperidin-1-yl)azetidin-1-

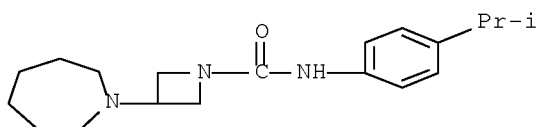
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RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of azetidine ring compds. as EDG-5 antagonists for treatment of hypertension, myocardial infarction, etc.)

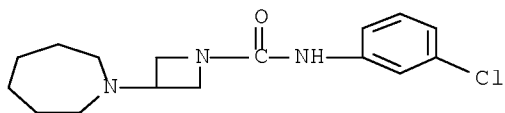
RN 858653-43-7 CAPLUS

CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-[4-(1-methylethyl)phenyl]- (CA INDEX NAME)



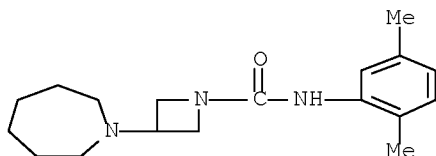
RN 858653-44-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-(hexahydro-1H-azepin-1-yl)- (CA INDEX NAME)



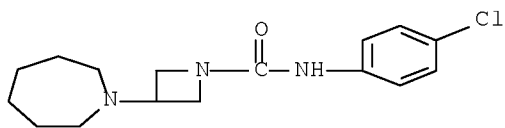
RN 858653-45-9 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,5-dimethylphenyl)-3-(hexahydro-1H-azepin-1-yl)- (CA INDEX NAME)



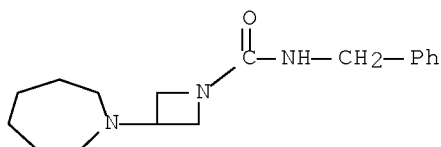
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CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-(hexahydro-1H-azepin-1-yl)- (CA INDEX NAME)



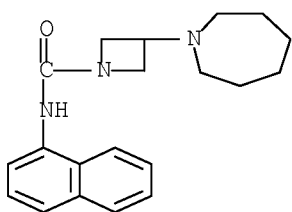
RN 858653-47-1 CAPLUS

CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-(phenylmethyl)- (CA INDEX NAME)



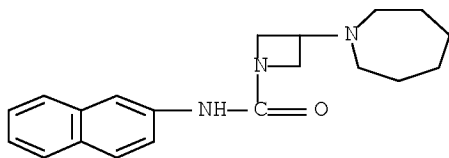
RN 858653-48-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-1-naphthalenyl- (CA INDEX NAME)



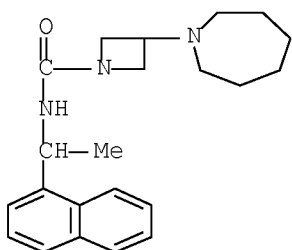
RN 858653-49-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-2-naphthalenyl-
(CA INDEX NAME)



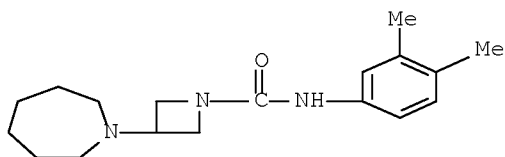
RN 858653-50-6 CAPLUS

CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-[1-(1-naphthalenyl)ethyl]- (CA INDEX NAME)



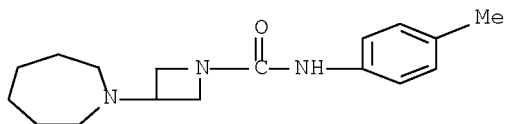
RN 858653-51-7 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,4-dimethylphenyl)-3-(hexahydro-1H-azepin-1-yl)- (CA INDEX NAME)



RN 858653-52-8 CAPLUS

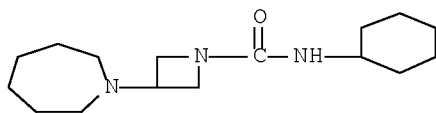
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RN 858653-53-9 CAPLUS

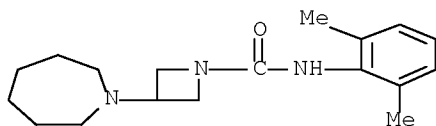
CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(hexahydro-1H-azepin-1-yl)- (CA

INDEX NAME)



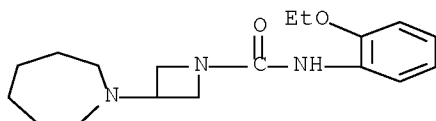
RN 858653-54-0 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-(hexahydro-1H-azepin-1-yl)-
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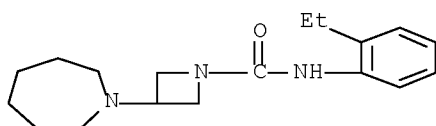
RN 858653-55-1 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-ethoxyphenyl)-3-(hexahydro-1H-azepin-1-yl)-
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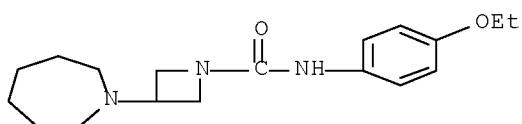
RN 858653-56-2 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-ethylphenyl)-3-(hexahydro-1H-azepin-1-yl)-
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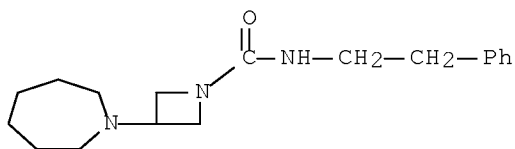


RN 858653-57-3 CAPLUS

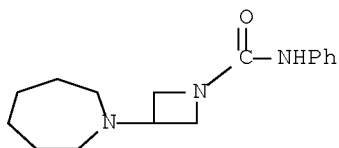
CN 1-Azetidinecarboxamide, N-(4-ethoxyphenyl)-3-(hexahydro-1H-azepin-1-yl)-
(CA INDEX NAME)



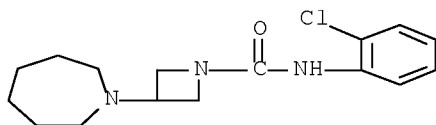
RN 858653-58-4 CAPLUS
CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-(2-phenylethyl)-
(CA INDEX NAME)



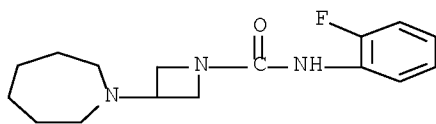
RN 858653-59-5 CAPLUS
CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-phenyl- (CA INDEX
NAME)



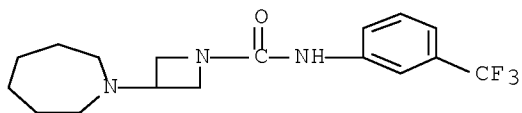
RN 858653-60-8 CAPLUS
CN 1-Azetidinecarboxamide, N-(2-chlorophenyl)-3-(hexahydro-1H-azepin-1-yl)-
(CA INDEX NAME)



RN 858653-61-9 CAPLUS
CN 1-Azetidinecarboxamide, N-(2-fluorophenyl)-3-(hexahydro-1H-azepin-1-yl)-
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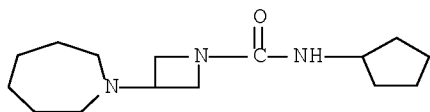


RN 858653-62-0 CAPLUS
CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



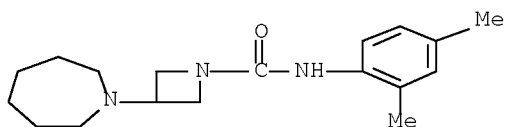
RN 858653-63-1 CAPLUS

CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(hexahydro-1H-azepin-1-yl)- 4-(trifluoromethyl)- (CA INDEX NAME)



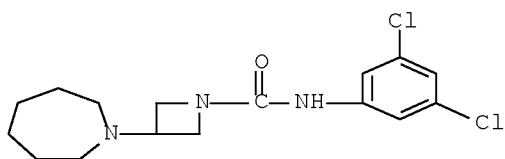
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CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(hexahydro-1H-azepin-1-yl)- (CA INDEX NAME)



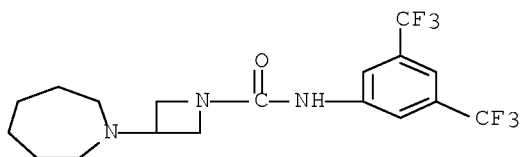
RN 858653-65-3 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(hexahydro-1H-azepin-1-yl)- (CA INDEX NAME)



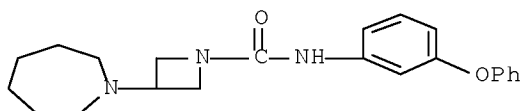
RN 858653-66-4 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(hexahydro-1H-azepin-1-yl)- (CA INDEX NAME)



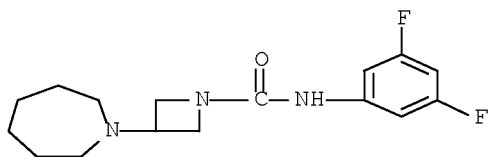
RN 858653-67-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-(3-phenoxyphenyl)-
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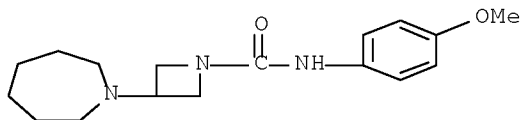
RN 858653-68-6 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-(hexahydro-1H-azepin-1-yl)-
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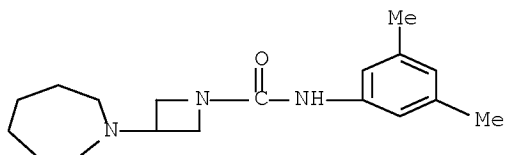
RN 858653-69-7 CAPLUS

CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-(4-methoxyphenyl)-
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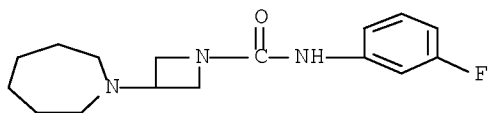
RN 858653-70-0 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-(hexahydro-1H-azepin-1-yl)-
(CA INDEX NAME)



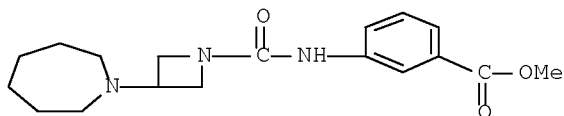
RN 858653-71-1 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-fluorophenyl)-3-(hexahydro-1H-azepin-1-yl)-
(CA INDEX NAME)



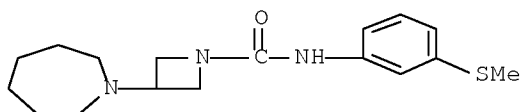
RN 858653-72-2 CAPLUS

CN Benzoic acid, 3-[[[3-(hexahydro-1H-azepin-1-yl)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



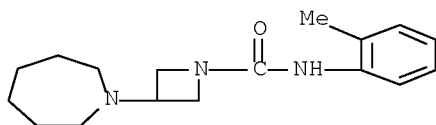
RN 858653-73-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-[3-(methylthio)phenyl]- (CA INDEX NAME)



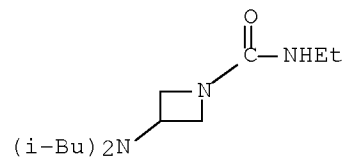
RN 858653-74-4 CAPLUS

CN 1-Azetidinecarboxamide, 3-(hexahydro-1H-azepin-1-yl)-N-(2-methylphenyl)- (CA INDEX NAME)



RN 858653-75-5 CAPLUS

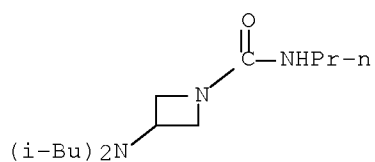
CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-ethyl- (CA INDEX NAME)



RN 858653-76-6 CAPLUS

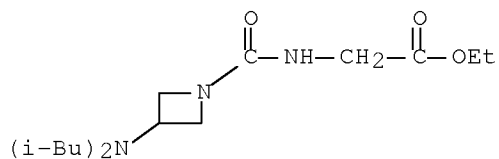
CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-propyl- (CA INDEX NAME)

NAME)



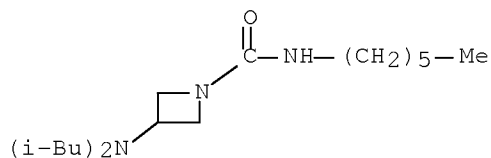
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CN Glycine, N-[[3-[bis(2-methylpropyl)amino]-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)



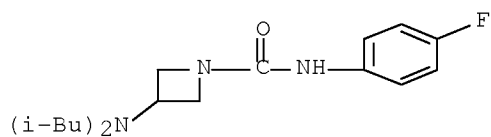
RN 858653-78-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-hexyl- (CA INDEX NAME)



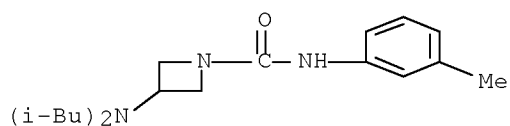
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CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-(4-fluorophenyl)- (CA INDEX NAME)



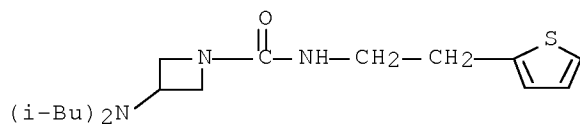
RN 858653-80-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-(3-methylphenyl)- (CA INDEX NAME)



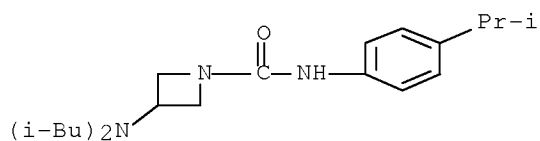
RN 858653-81-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-[2-(2-thienyl)ethyl]- (CA INDEX NAME)



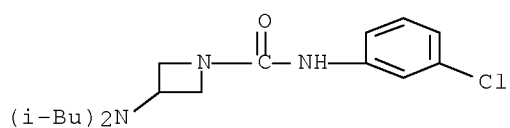
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CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-[4-(1-methylethyl)phenyl]- (CA INDEX NAME)



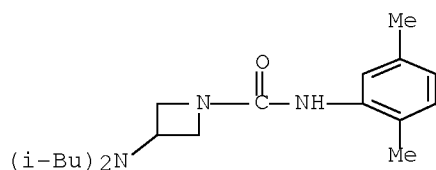
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CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-(3-chlorophenyl)- (CA INDEX NAME)



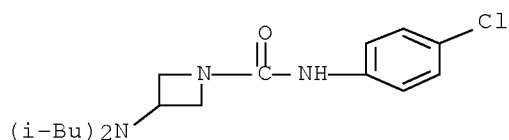
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CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-(2,5-dimethylphenyl)- (CA INDEX NAME)



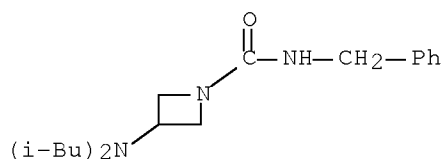
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CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-(4-chlorophenyl)-
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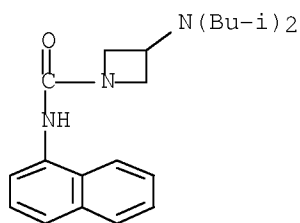
RN 858653-86-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-(phenylmethyl)-
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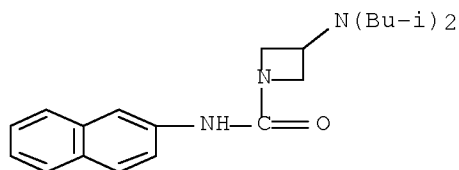
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CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-1-naphthalenyl-
(CA INDEX NAME)



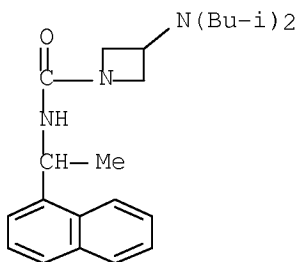
RN 858653-88-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-2-naphthalenyl-
(CA INDEX NAME)



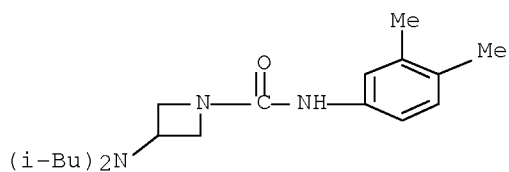
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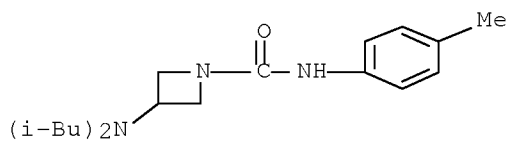
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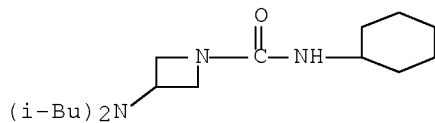
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CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-(4-methylphenyl)- (CA INDEX NAME)



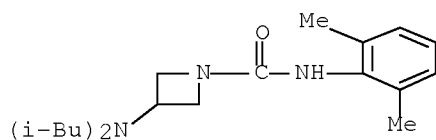
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CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-cyclohexyl- (CA INDEX NAME)



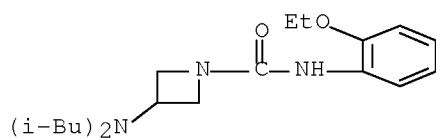
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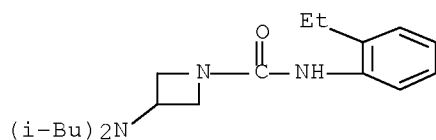
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CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-(2-ethoxyphenyl)-
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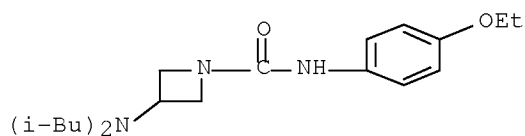
RN 858653-95-9 CAPLUS

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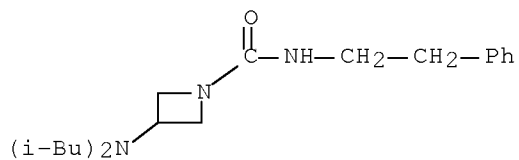
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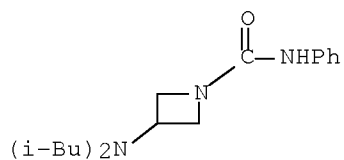


RN 858653-97-1 CAPLUS

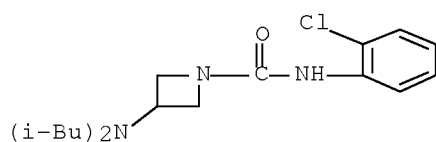
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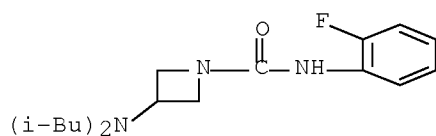
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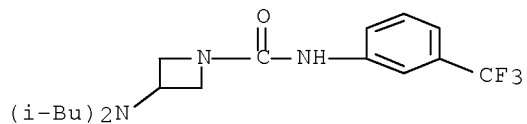
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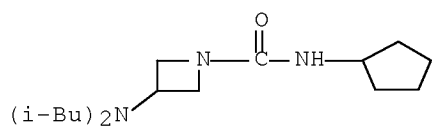
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RN 858654-01-0 CAPLUS
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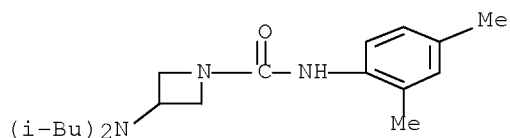


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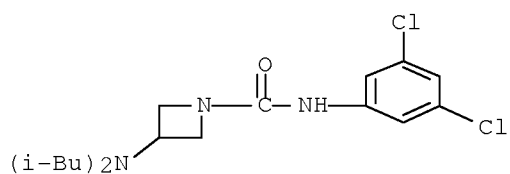
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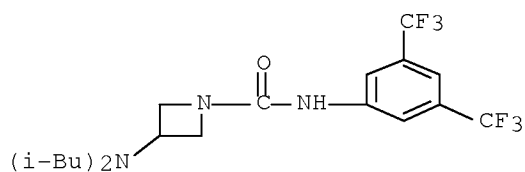
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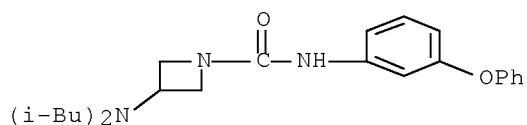
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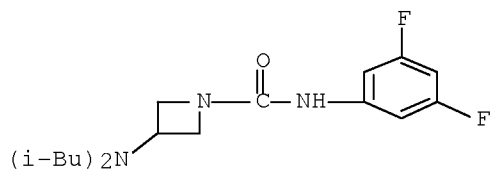


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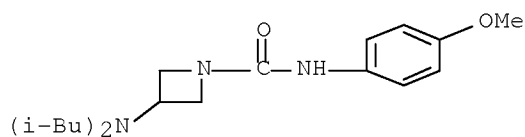
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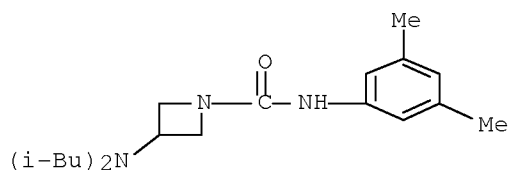
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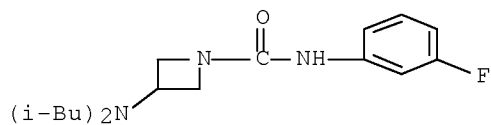
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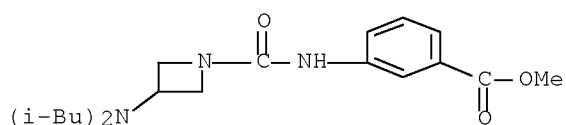
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CN 1-Azetidinecarboxamide, 3-[bis(2-methylpropyl)amino]-N-(3,5-dimethylphenyl)- (CA INDEX NAME)



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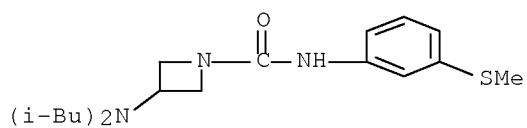


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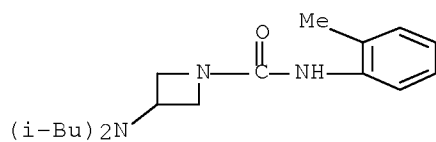
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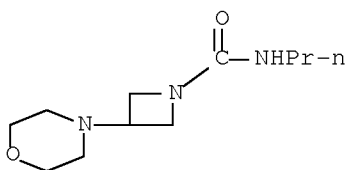
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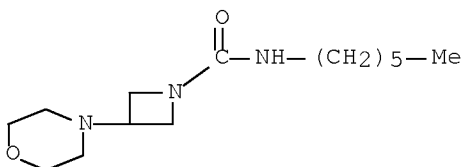
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CN 1-Azetidinecarboxamide, 3-(4-morpholinyl)-N-propyl- (CA INDEX NAME)

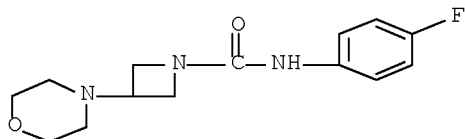


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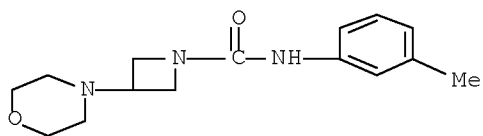
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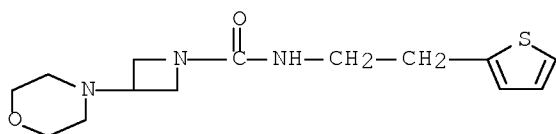
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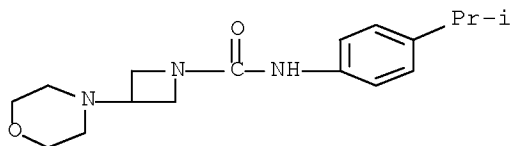
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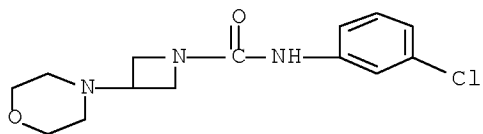
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CN 1-Azetidinecarboxamide, 3-(4-morpholinyl)-N-[2-(2-thienyl)ethyl]- (CA INDEX NAME)



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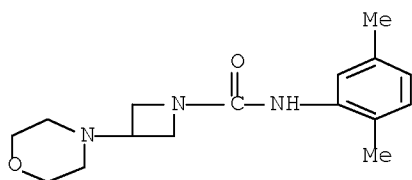


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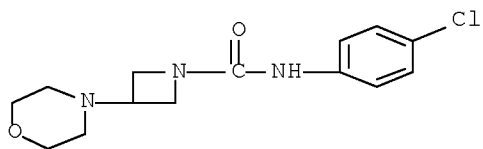
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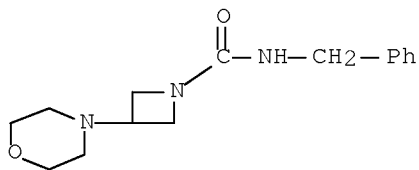
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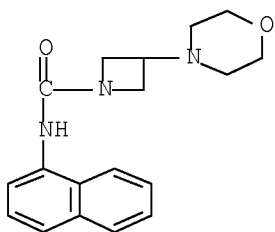
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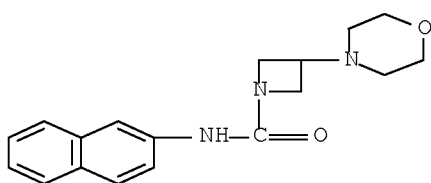
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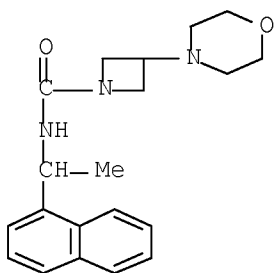
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CN 1-Azetidinecarboxamide, 3-(4-morpholinyl)-N-2-naphthalenyl- (CA INDEX NAME)



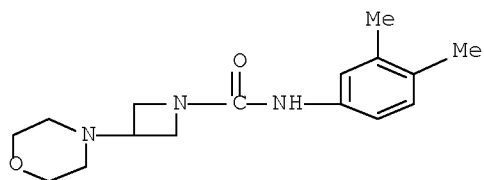
RN 858654-27-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-(4-morpholinyl)-N-[1-(1-naphthalenyl)ethyl]- (CA INDEX NAME)



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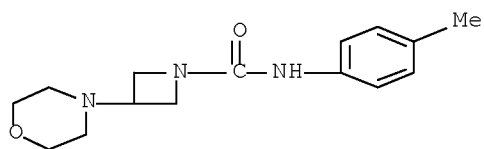
CN 1-Azetidinecarboxamide, N-(3,4-dimethylphenyl)-3-(4-morpholinyl)- (CA INDEX NAME)



RN 858654-30-5 CAPLUS

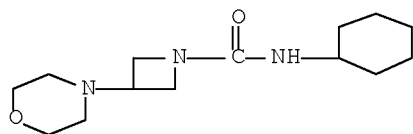
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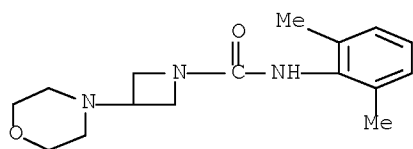
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CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(4-morpholinyl)- (CA INDEX NAME)



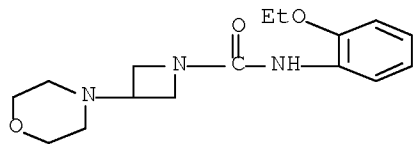
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CN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-(4-morpholinyl)- (CA INDEX NAME)



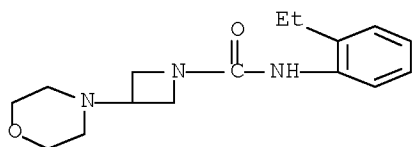
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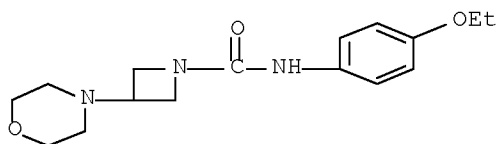


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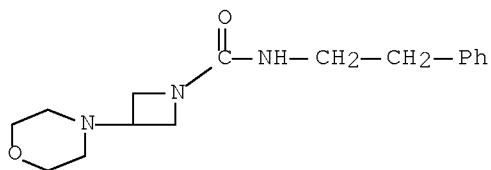
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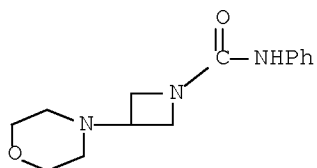
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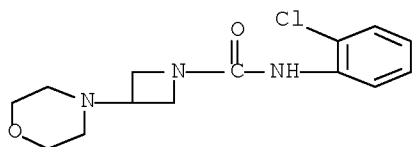
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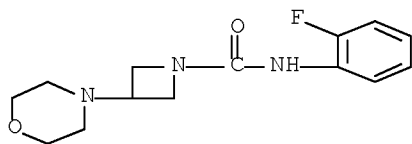


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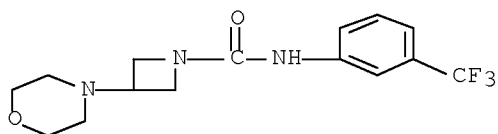
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CN 1-Azetidinecarboxamide, N-(2-fluorophenyl)-3-(4-morpholinyl)- (CA INDEX NAME)



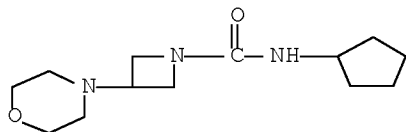
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CN 1-Azetidinecarboxamide, 3-(4-morpholinyl)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



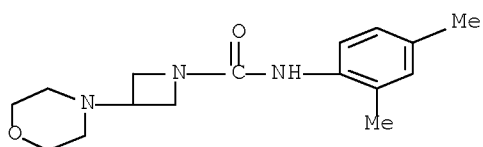
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CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(4-morpholinyl)- (CA INDEX NAME)



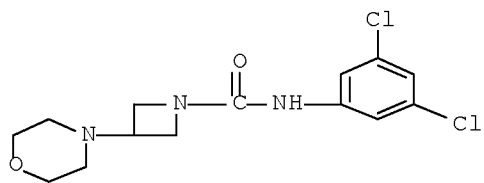
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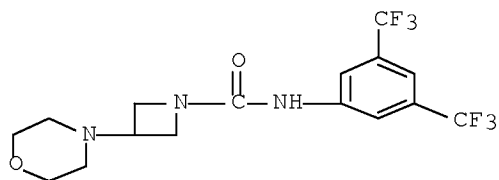
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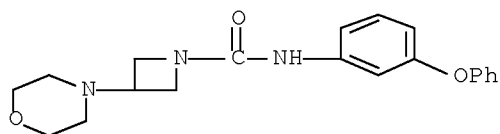
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(4-morpholinyl)- (CA INDEX NAME)



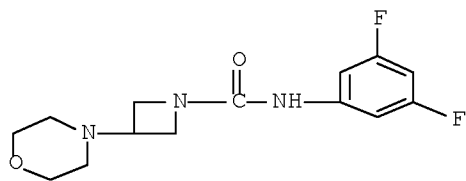
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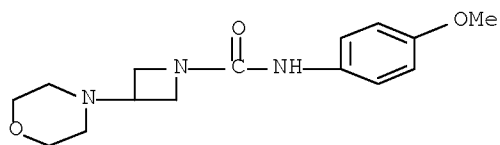
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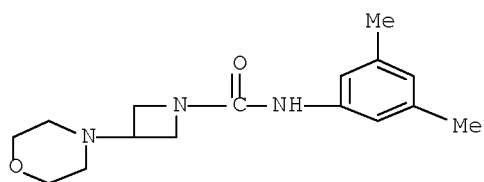
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NAME)



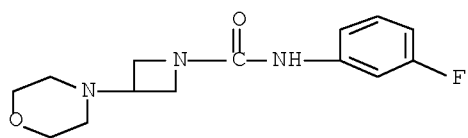
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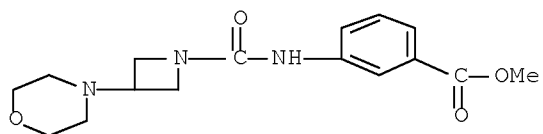
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CN 1-Azetidinecarboxamide, N-(3-fluorophenyl)-3-(4-morpholinyl)- (CA INDEX NAME)



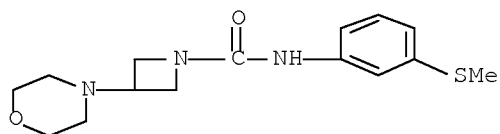
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CN Benzoic acid, 3-[[[3-(4-morpholinyl)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



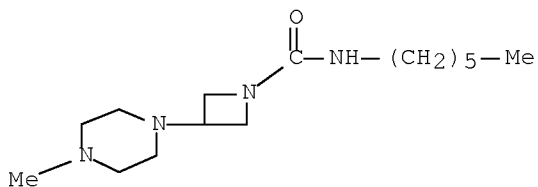
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CN 1-Azetidinecarboxamide, N-[3-(methylthio)phenyl]-3-(4-morpholinyl)- (CA INDEX NAME)



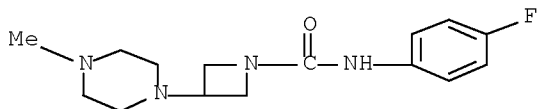
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CN 1-Azetidinecarboxamide, N-hexyl-3-(4-methyl-1-piperazinyl)- (CA INDEX NAME)



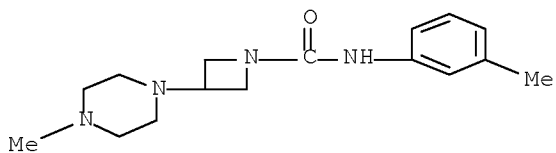
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CN 1-Azetidinecarboxamide, N-(4-fluorophenyl)-3-(4-methyl-1-piperazinyl)- (CA INDEX NAME)



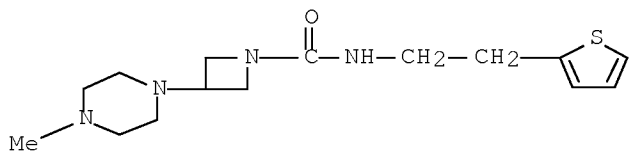
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CN 1-Azetidinecarboxamide, N-(3-methylphenyl)-3-(4-methyl-1-piperazinyl)- (CA INDEX NAME)



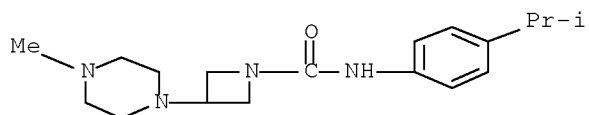
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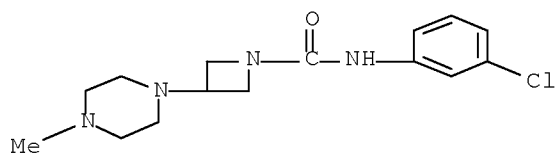
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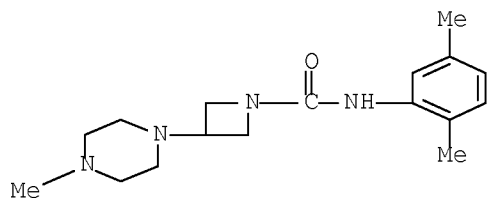
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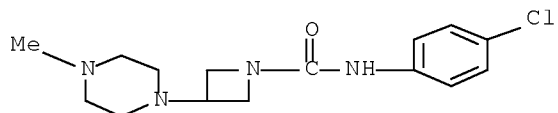
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CN 1-Azetidinecarboxamide, N-(2,5-dimethylphenyl)-3-(4-methyl-1-piperazinyl)- (CA INDEX NAME)



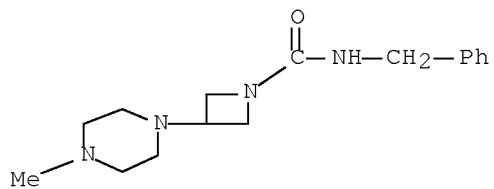
RN 858654-59-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-(4-methyl-1-piperazinyl)- (CA INDEX NAME)



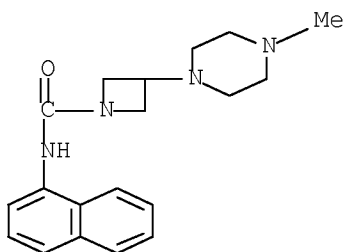
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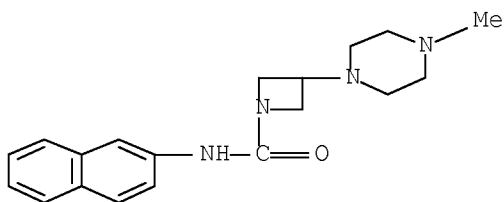
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CN 1-Azetidinecarboxamide, 3-(4-methyl-1-piperazinyl)-N-1-naphthalenyl- (CA INDEX NAME)



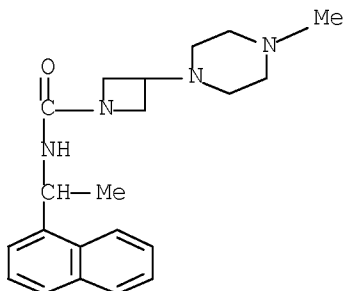
RN 858654-62-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(4-methyl-1-piperazinyl)-N-2-naphthalenyl- (CA INDEX NAME)



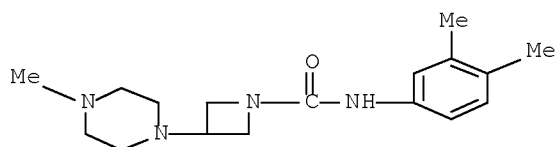
RN 858654-63-4 CAPLUS

CN 1-Azetidinecarboxamide, 3-(4-methyl-1-piperazinyl)-N-[1-(1-naphthalenyl)ethyl]- (CA INDEX NAME)



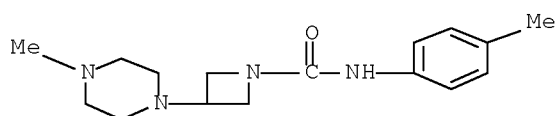
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CN 1-Azetidinecarboxamide, N-(3,4-dimethylphenyl)-3-(4-methyl-1-piperazinyl)-
(CA INDEX NAME)



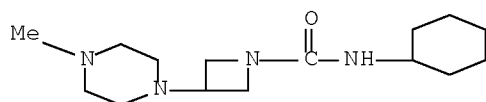
RN 858654-65-6 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-methylphenyl)-3-(4-methyl-1-piperazinyl)-
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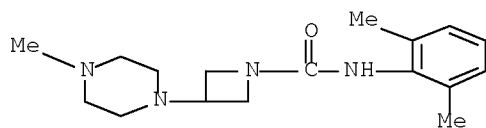
RN 858654-66-7 CAPLUS

CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(4-methyl-1-piperazinyl)- (CA
INDEX NAME)



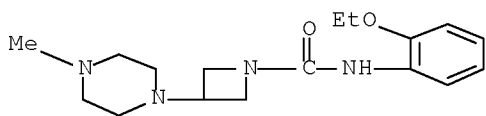
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CN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-(4-methyl-1-piperazinyl)-
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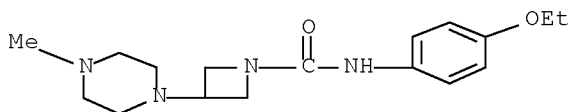
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CN 1-Azetidinecarboxamide, N-(2-ethoxyphenyl)-3-(4-methyl-1-piperazinyl)-
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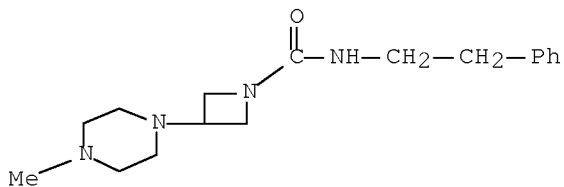
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CN 1-Azetidinecarboxamide, N-(4-ethoxyphenyl)-3-(4-methyl-1-piperazinyl)-
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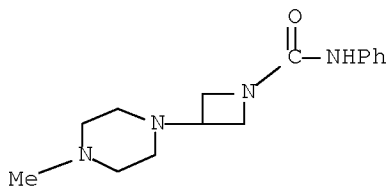
RN 858654-70-3 CAPLUS

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INDEX NAME)



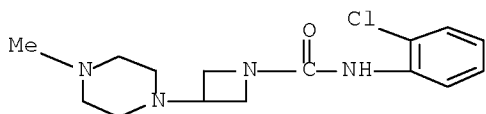
RN 858654-71-4 CAPLUS

CN 1-Azetidinecarboxamide, 3-(4-methyl-1-piperazinyl)-N-phenyl- (CA INDEX
NAME)



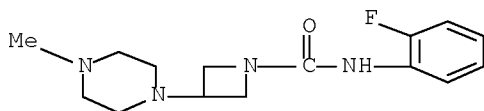
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CN 1-Azetidinecarboxamide, N-(2-chlorophenyl)-3-(4-methyl-1-piperazinyl)-
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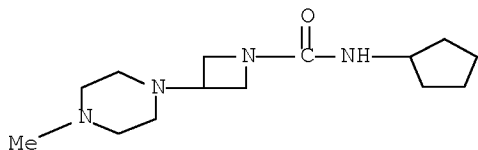
RN 858654-73-6 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-fluorophenyl)-3-(4-methyl-1-piperazinyl)-
(CA INDEX NAME)



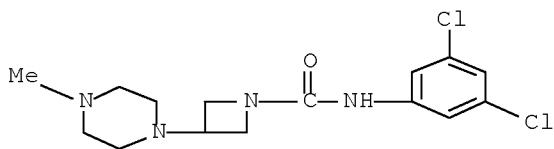
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INDEX NAME)



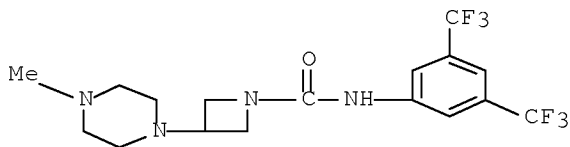
RN 858654-75-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(4-methyl-1-piperazinyl)-
(CA INDEX NAME)



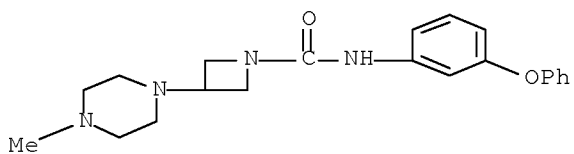
RN 858654-76-9 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(4-methyl-1-
piperazinyl)- (CA INDEX NAME)



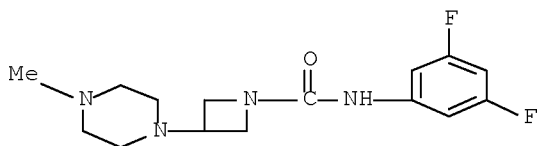
RN 858654-77-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-(4-methyl-1-piperazinyl)-N-(3-phenoxyphenyl)-
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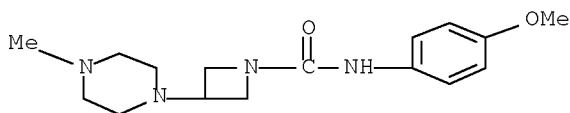
RN 858654-78-1 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-(4-methyl-1-piperazinyl)-
(CA INDEX NAME)



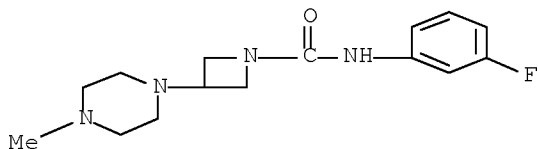
RN 858654-79-2 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-methoxyphenyl)-3-(4-methyl-1-piperazinyl)-
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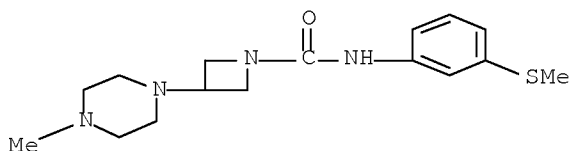
RN 858654-80-5 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-fluorophenyl)-3-(4-methyl-1-piperazinyl)-
(CA INDEX NAME)

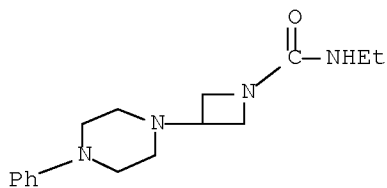


RN 858654-81-6 CAPLUS

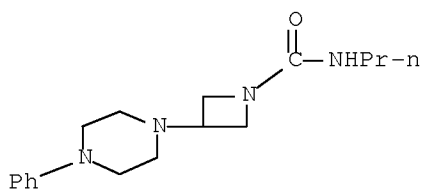
CN 1-Azetidinecarboxamide, 3-(4-methyl-1-piperazinyl)-N-[3-(methylthio)phenyl]-
(CA INDEX NAME)



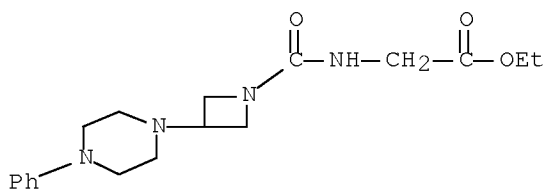
RN 858654-82-7 CAPLUS
 CN 1-Azetidinecarboxamide, N-ethyl-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)



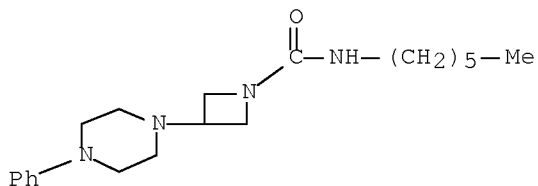
RN 858654-83-8 CAPLUS
 CN 1-Azetidinecarboxamide, 3-(4-phenyl-1-piperazinyl)-N-propyl- (CA INDEX NAME)



RN 858654-84-9 CAPLUS
 CN Glycine, N-[[3-(4-phenyl-1-piperazinyl)-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)

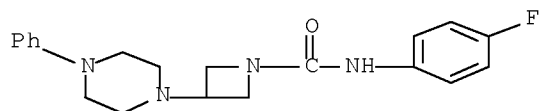


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 CN 1-Azetidinecarboxamide, N-hexyl-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)



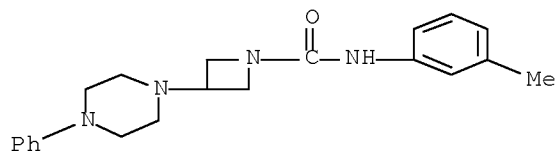
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CN 1-Azetidinecarboxamide, N-(4-fluorophenyl)-3-(4-phenyl-1-piperazinyl)-
(CA INDEX NAME)



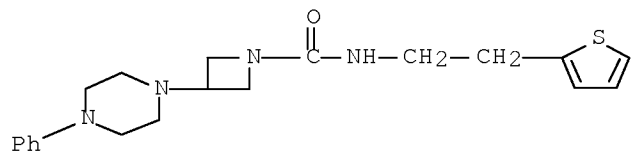
RN 858654-87-2 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-methylphenyl)-3-(4-phenyl-1-piperazinyl)-
(CA INDEX NAME)



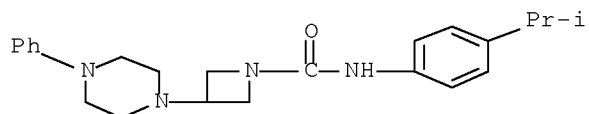
RN 858654-88-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(4-phenyl-1-piperazinyl)-N-[2-(2-thienyl)ethyl]-
(CA INDEX NAME)



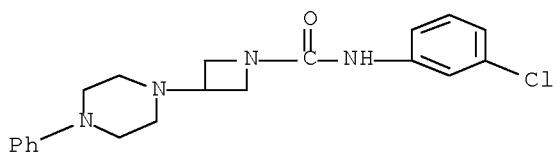
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CN 1-Azetidinecarboxamide, N-[4-(1-methylethyl)phenyl]-3-(4-phenyl-1-piperazinyl)-
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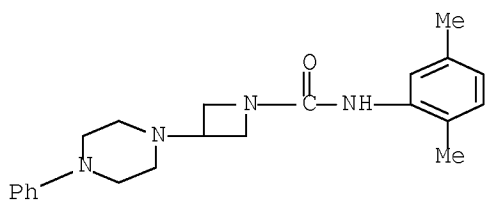
RN 858654-90-7 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-(4-phenyl-1-piperazinyl)-
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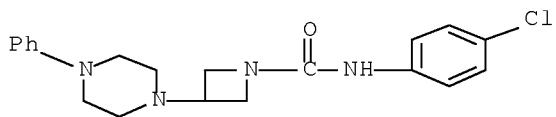
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CN 1-Azetidinecarboxamide, N-(2,5-dimethylphenyl)-3-(4-phenyl-1-piperazinyl)-
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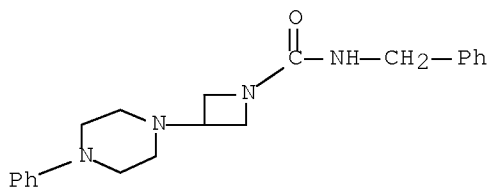
RN 858654-92-9 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-(4-phenyl-1-piperazinyl)-
(CA INDEX NAME)



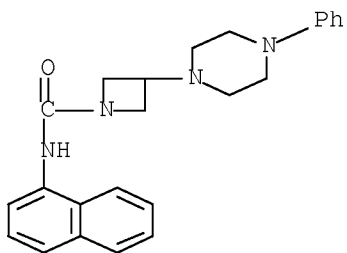
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INDEX NAME)



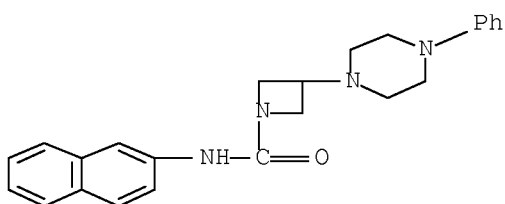
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CN 1-Azetidinecarboxamide, N-1-naphthalenyl-3-(4-phenyl-1-piperazinyl)- (CA
INDEX NAME)



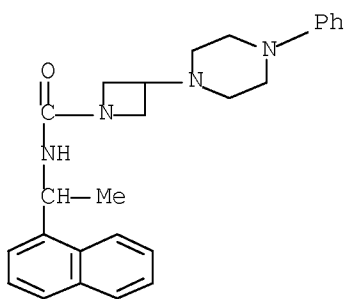
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CN 1-Azetidinecarboxamide, N-2-naphthalenyl-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)



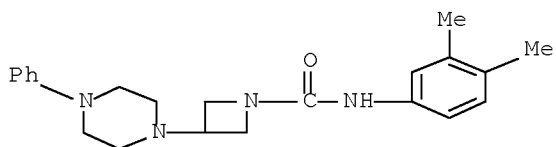
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CN 1-Azetidinecarboxamide, N-[1-(1-naphthalenyl)ethyl]-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)



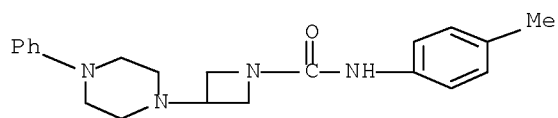
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CN 1-Azetidinecarboxamide, N-(3,4-dimethylphenyl)-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)



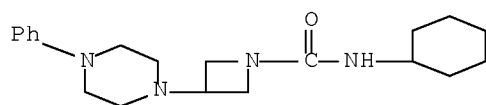
RN 858655-00-2 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-methylphenyl)-3-(4-phenyl-1-piperazinyl)-
(CA INDEX NAME)



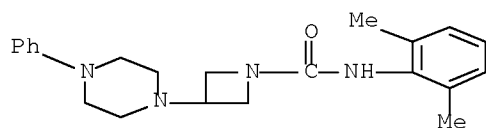
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INDEX NAME)



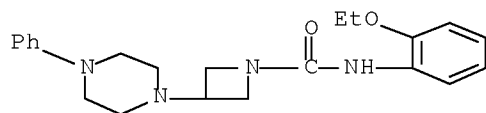
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CN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-(4-phenyl-1-piperazinyl)-
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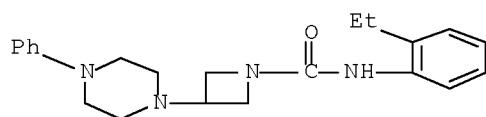
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CN 1-Azetidinecarboxamide, N-(2-ethoxyphenyl)-3-(4-phenyl-1-piperazinyl)-
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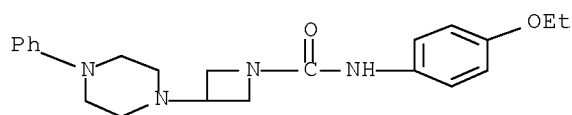
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INDEX NAME)



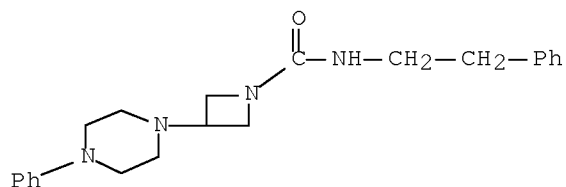
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CN 1-Azetidinecarboxamide, N-(4-ethoxyphenyl)-3-(4-phenyl-1-piperazinyl)-
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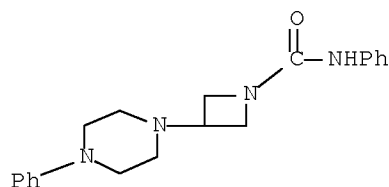
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CN 1-Azetidinecarboxamide, N-(2-phenylethyl)-3-(4-phenyl-1-piperazinyl)- (CA
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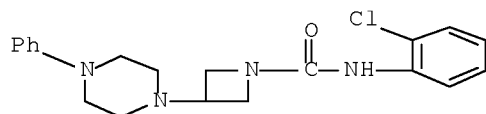
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CN 1-Azetidinecarboxamide, N-phenyl-3-(4-phenyl-1-piperazinyl)- (CA INDEX
NAME)



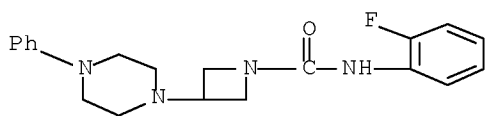
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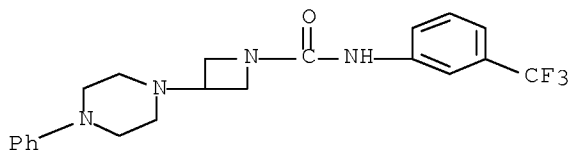
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CN 1-Azetidinecarboxamide, N-(2-fluorophenyl)-3-(4-phenyl-1-piperazinyl)-
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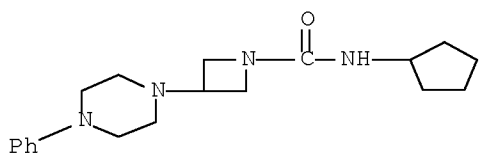
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CN 1-Azetidinecarboxamide, 3-(4-phenyl-1-piperazinyl)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



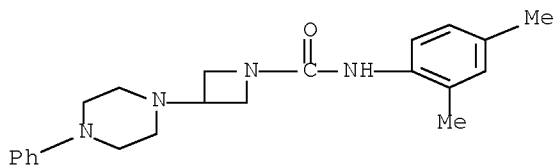
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CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)



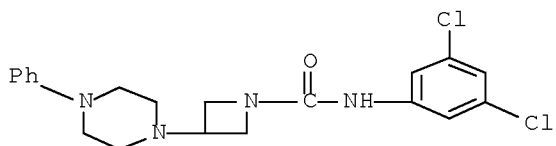
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CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)

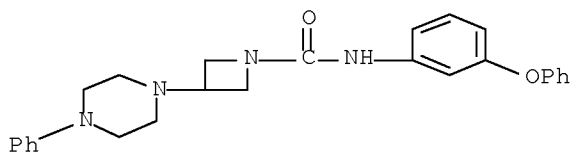


RN 858655-13-7 CAPLUS

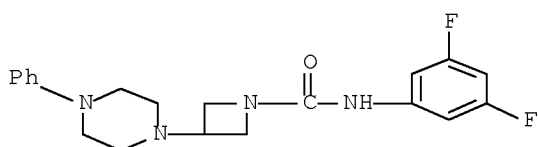
CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)



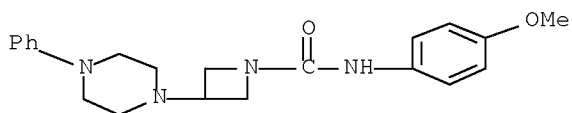
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CN 1-Azetidinecarboxamide, N-(3-phenoxyphenyl)-3-(4-phenyl-1-piperazinyl)-
(CA INDEX NAME)



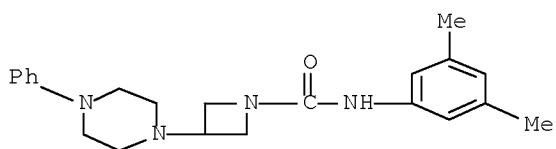
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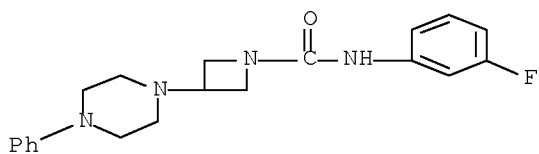
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CN 1-Azetidinecarboxamide, N-(4-methoxyphenyl)-3-(4-phenyl-1-piperazinyl)-
(CA INDEX NAME)



RN 858655-17-1 CAPLUS
CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-(4-phenyl-1-piperazinyl)-
(CA INDEX NAME)

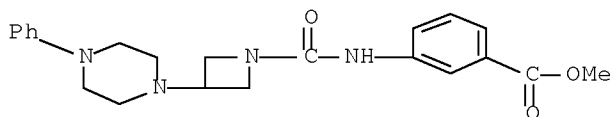


RN 858655-18-2 CAPLUS
CN 1-Azetidinecarboxamide, N-(3-fluorophenyl)-3-(4-phenyl-1-piperazinyl)-
(CA INDEX NAME)



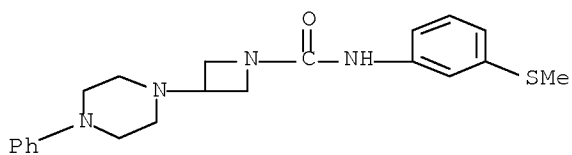
RN 858655-19-3 CAPLUS

CN Benzoic acid, 3-[[[3-(4-phenyl-1-piperazinyl)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



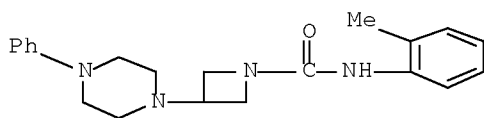
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CN 1-Azetidinecarboxamide, N-[3-(methylthio)phenyl]-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)



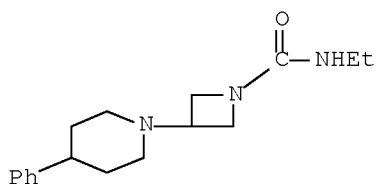
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CN 1-Azetidinecarboxamide, N-(2-methylphenyl)-3-(4-phenyl-1-piperazinyl)- (CA INDEX NAME)

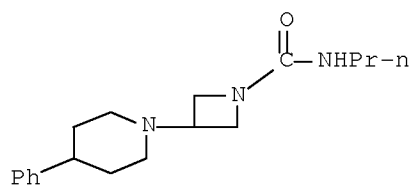


RN 858655-22-8 CAPLUS

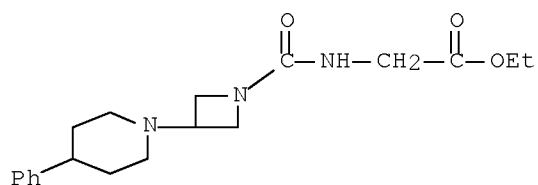
CN 1-Azetidinecarboxamide, N-ethyl-3-(4-phenyl-1-piperidinyl)- (CA INDEX NAME)



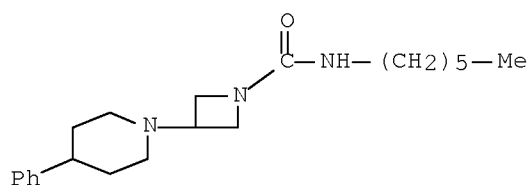
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 CN 1-Azetidinecarboxamide, 3-(4-phenyl-1-piperidiny1)-N-propyl- (CA INDEX NAME)



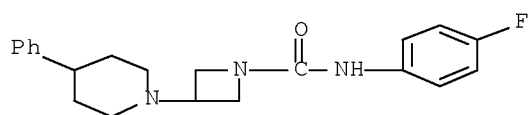
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 CN Glycine, N-[[3-(4-phenyl-1-piperidiny1)-1-azetidiny1]carbonyl]-, ethyl ester (CA INDEX NAME)



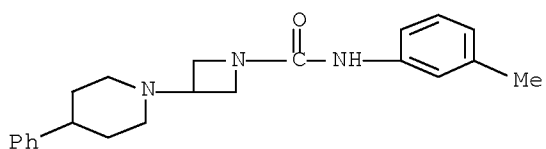
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 CN 1-Azetidinecarboxamide, N-hexyl-3-(4-phenyl-1-piperidiny1)- (CA INDEX NAME)



RN 858655-26-2 CAPLUS
 CN 1-Azetidinecarboxamide, N-(4-fluorophenyl)-3-(4-phenyl-1-piperidiny1)- (CA INDEX NAME)

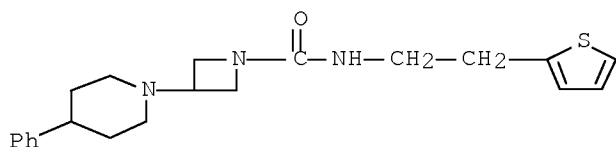


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 CN 1-Azetidinecarboxamide, N-(3-methylphenyl)-3-(4-phenyl-1-piperidiny1)- (CA INDEX NAME)



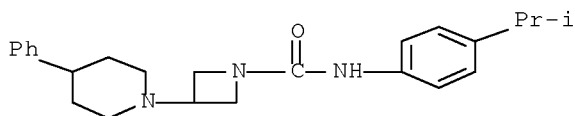
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CN 1-Azetidinecarboxamide, 3-(4-phenyl-1-piperidinyl)-N-[2-(2-thienyl)ethyl]-
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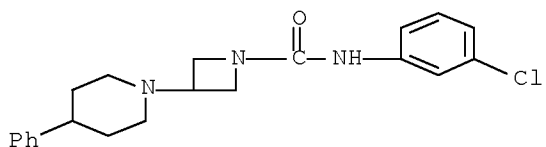
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CN 1-Azetidinecarboxamide, N-[4-(1-methylethyl)phenyl]-3-(4-phenyl-1-piperidinyl)-
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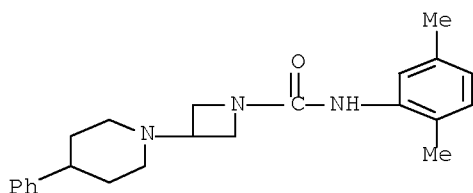
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CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-(4-phenyl-1-piperidinyl)-
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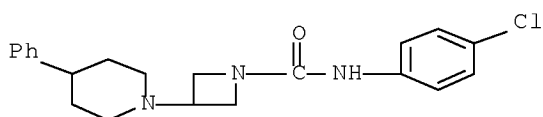
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CN 1-Azetidinecarboxamide, N-(2,5-dimethylphenyl)-3-(4-phenyl-1-piperidinyl)-
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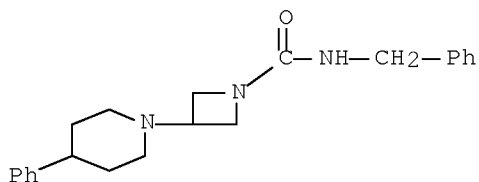
RN 858655-32-0 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-(4-phenyl-1-piperidinyl)-
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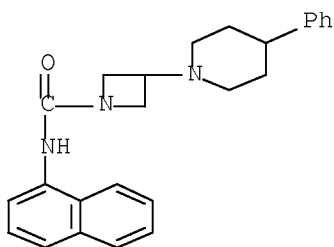
RN 858655-33-1 CAPLUS

CN 1-Azetidinecarboxamide, N-(phenylmethyl)-3-(4-phenyl-1-piperidinyl)- (CA
INDEX NAME)



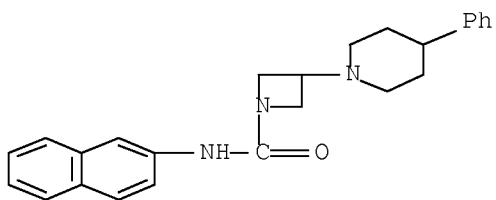
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CN 1-Azetidinecarboxamide, N-1-naphthalenyl-3-(4-phenyl-1-piperidinyl)- (CA
INDEX NAME)



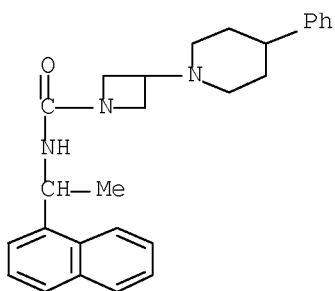
RN 858655-35-3 CAPLUS

CN 1-Azetidinecarboxamide, N-2-naphthalenyl-3-(4-phenyl-1-piperidinyl)- (CA
INDEX NAME)



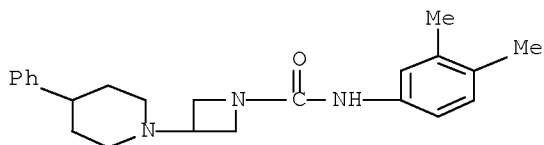
RN 858655-36-4 CAPLUS

CN 1-Azetidinecarboxamide, N-[1-(1-naphthalenyl)ethyl]-3-(4-phenyl-1-piperidiny)- (CA INDEX NAME)



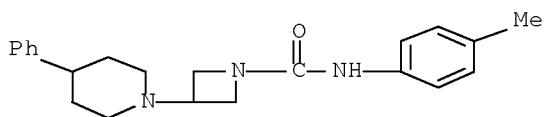
RN 858655-37-5 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,4-dimethylphenyl)-3-(4-phenyl-1-piperidiny)- (CA INDEX NAME)



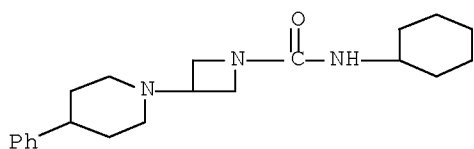
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CN 1-Azetidinecarboxamide, N-(4-methylphenyl)-3-(4-phenyl-1-piperidiny)- (CA INDEX NAME)



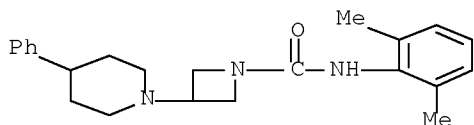
RN 858655-39-7 CAPLUS

CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(4-phenyl-1-piperidiny)- (CA INDEX NAME)



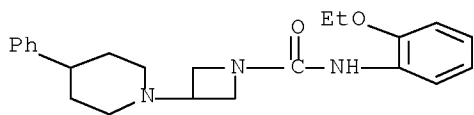
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CN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-(4-phenyl-1-piperidinyl)-
(CA INDEX NAME)



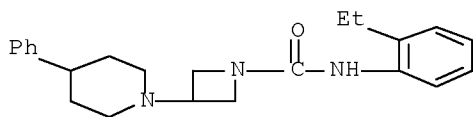
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CN 1-Azetidinecarboxamide, N-(2-ethoxyphenyl)-3-(4-phenyl-1-piperidinyl)-
(CA INDEX NAME)



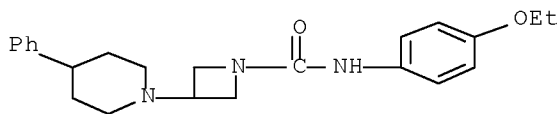
RN 858655-42-2 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-ethylphenyl)-3-(4-phenyl-1-piperidinyl)- (CA
INDEX NAME)



RN 858655-43-3 CAPLUS

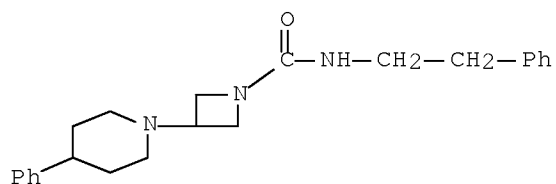
CN 1-Azetidinecarboxamide, N-(4-ethoxyphenyl)-3-(4-phenyl-1-piperidinyl)-
(CA INDEX NAME)



RN 858655-44-4 CAPLUS

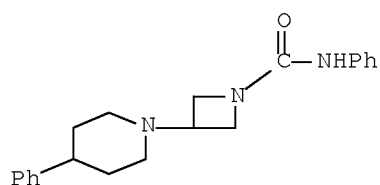
CN 1-Azetidinecarboxamide, N-(2-phenylethyl)-3-(4-phenyl-1-piperidinyl)- (CA

INDEX NAME)



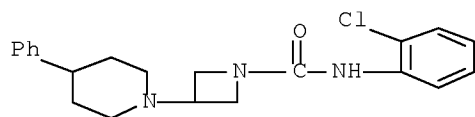
RN 858655-45-5 CAPLUS

CN 1-Azetidinecarboxamide, N-phenyl-3-(4-phenyl-1-piperidinyl)- (CA INDEX NAME)



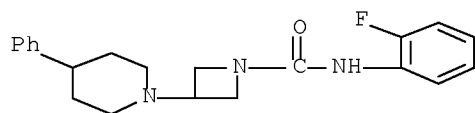
RN 858655-46-6 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-chlorophenyl)-3-(4-phenyl-1-piperidinyl)- (CA INDEX NAME)



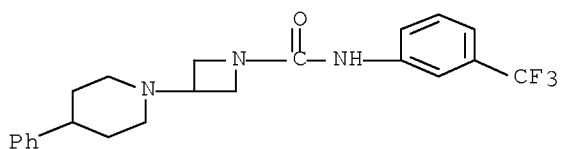
RN 858655-47-7 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-fluorophenyl)-3-(4-phenyl-1-piperidinyl)- (CA INDEX NAME)



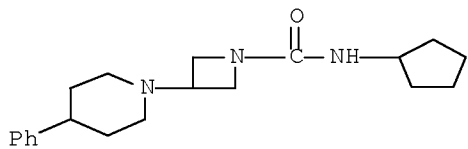
RN 858655-48-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-(4-phenyl-1-piperidinyl)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



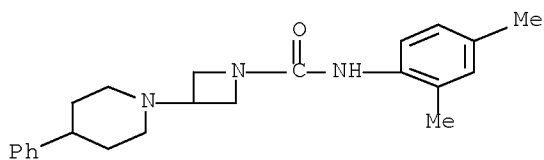
RN 858655-49-9 CAPLUS

CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(4-phenyl-1-piperidinyl)- (CA INDEX NAME)



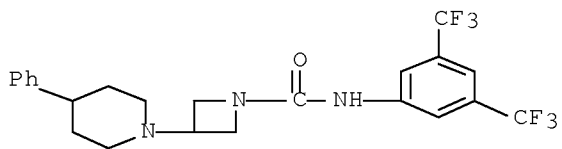
RN 858655-50-2 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-(4-phenyl-1-piperidinyl)- (CA INDEX NAME)



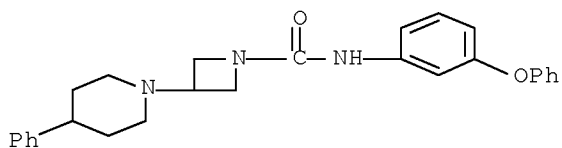
RN 858655-51-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(4-phenyl-1-piperidinyl)- (CA INDEX NAME)



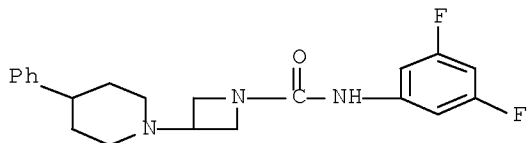
RN 858655-52-4 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-phenoxyphenyl)-3-(4-phenyl-1-piperidinyl)- (CA INDEX NAME)



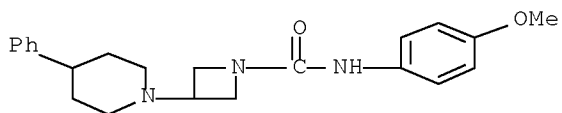
RN 858655-53-5 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-(4-phenyl-1-piperidiny)-
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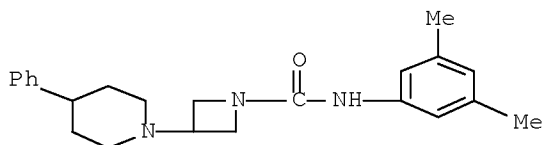
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CN 1-Azetidinecarboxamide, N-(4-methoxyphenyl)-3-(4-phenyl-1-piperidiny)-
(CA INDEX NAME)



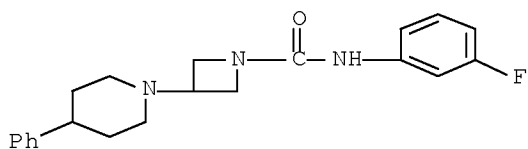
RN 858655-55-7 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-(4-phenyl-1-piperidiny)-
(CA INDEX NAME)



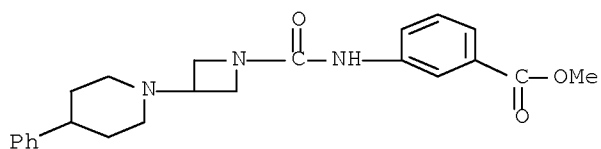
RN 858655-56-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-fluorophenyl)-3-(4-phenyl-1-piperidiny)-
(CA INDEX NAME)



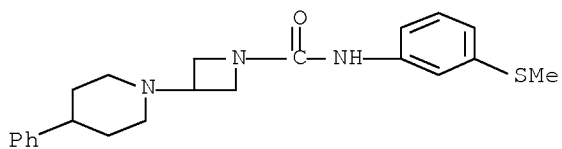
RN 858655-57-9 CAPLUS

CN Benzoic acid, 3-[[[3-(4-phenyl-1-piperidiny)-1-azetidiny]carbonyl]amino]-
, methyl ester (CA INDEX NAME)



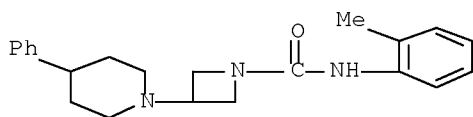
RN 858655-58-0 CAPLUS

CN 1-Azetidinecarboxamide, N-[3-(methylthio)phenyl]-3-(4-phenyl-1-piperidinyl)- (CA INDEX NAME)



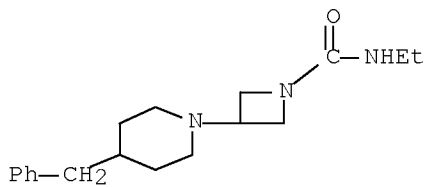
RN 858655-59-1 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-methylphenyl)-3-(4-phenyl-1-piperidinyl)- (CA INDEX NAME)



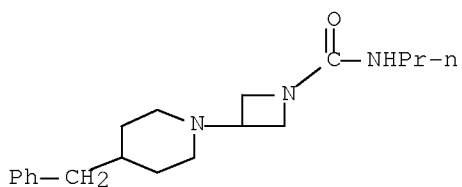
RN 858655-60-4 CAPLUS

CN 1-Azetidinecarboxamide, N-ethyl-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



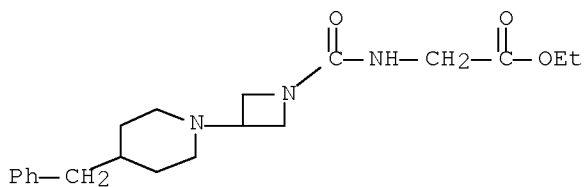
RN 858655-61-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-[4-(phenylmethyl)-1-piperidinyl]-N-propyl- (CA INDEX NAME)



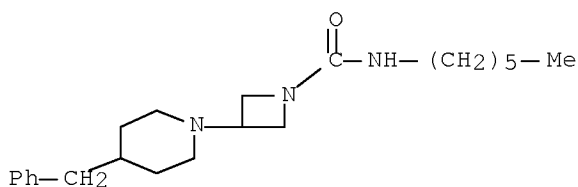
RN 858655-62-6 CAPLUS

CN Glycine, N-[3-[4-(phenylmethyl)-1-piperidinyl]-1-azetidiny]-, ethyl ester (9CI) (CA INDEX NAME)



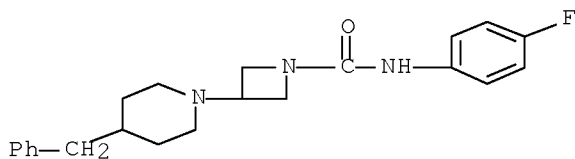
RN 858655-63-7 CAPLUS

CN 1-Azetidinecarboxamide, N-hexyl-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



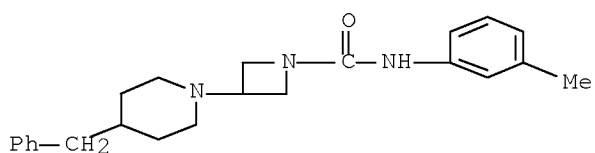
RN 858655-64-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(4-fluorophenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



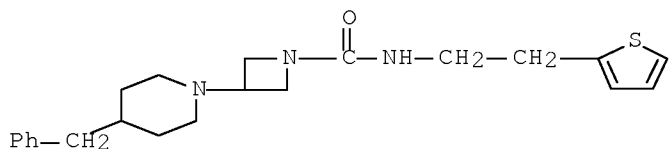
RN 858655-65-9 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-methylphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



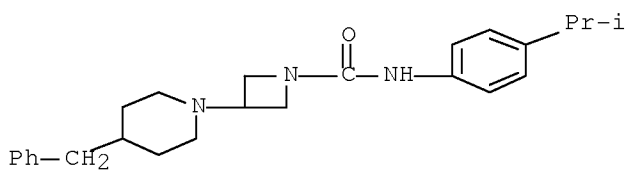
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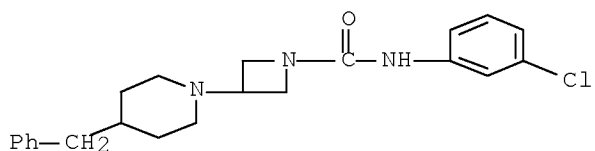
RN 858655-67-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[4-(1-methylethyl)phenyl]-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



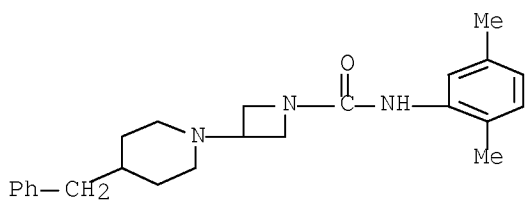
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CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



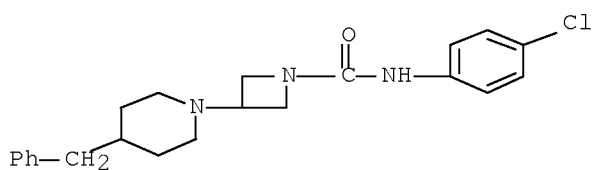
RN 858655-69-3 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,5-dimethylphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



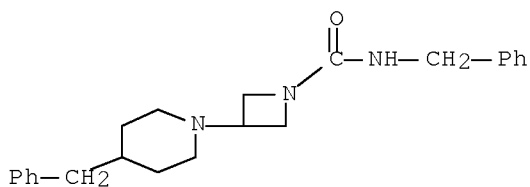
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CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



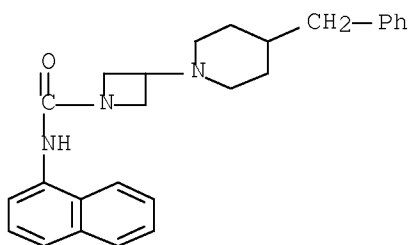
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CN 1-Azetidinecarboxamide, N-(phenylmethyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



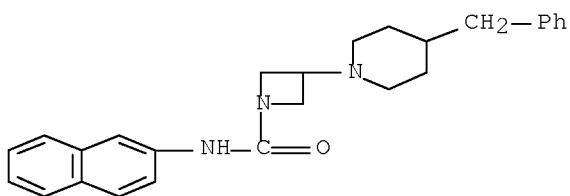
RN 858655-72-8 CAPLUS

CN 1-Azetidinecarboxamide, N-1-naphthalenyl-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



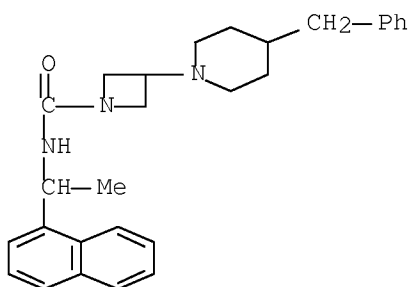
RN 858655-73-9 CAPLUS

CN 1-Azetidinecarboxamide, N-2-naphthalenyl-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



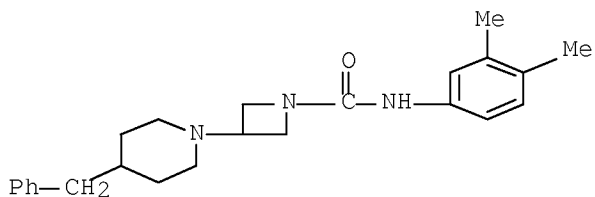
RN 858655-74-0 CAPLUS

CN 1-Azetidinecarboxamide, N-[1-(1-naphthalenyl)ethyl]-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



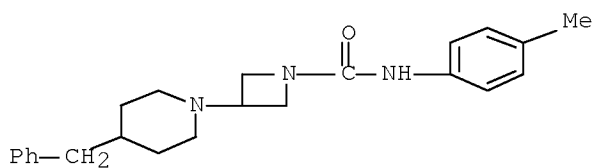
RN 858655-75-1 CAPLUS

CN 1-Azetidinecarboxamide, N-(3,4-dimethylphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



RN 858655-76-2 CAPLUS

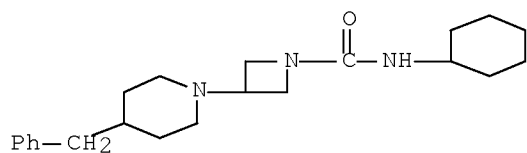
CN 1-Azetidinecarboxamide, N-(4-methylphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



RN 858655-77-3 CAPLUS

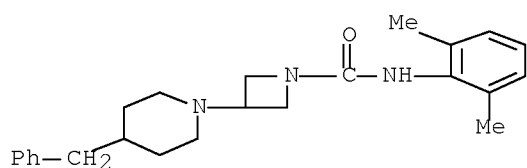
CN 1-Azetidinecarboxamide, N-cyclohexyl-3-[4-(phenylmethyl)-1-piperidinyl]-

(CA INDEX NAME)



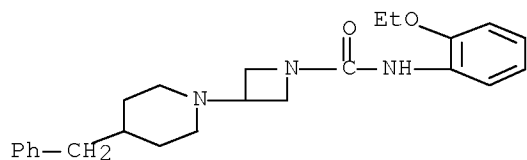
RN 858655-78-4 CAPLUS

CN 1-Azetidinecarboxamide, N-(2,6-dimethylphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



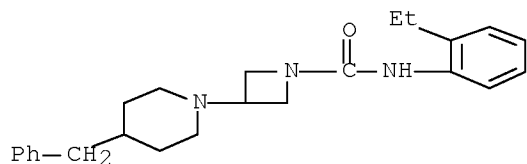
RN 858655-79-5 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-ethoxyphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



RN 858655-80-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-ethylphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



IT 858655-81-9P, 3-(4-Benzylpiperidin-1-yl)-N-(4-ethoxyphenyl)azetidine-1-carboxamide 858655-82-0P, 3-(4-Benzylpiperidin-1-yl)-N-(2-phenylethyl)azetidine-1-carboxamide 858655-83-1P, 3-(4-Benzylpiperidin-1-yl)-N-phenylazetidine-1-carboxamide 858655-84-2P, 3-(4-Benzylpiperidin-1-yl)-N-(2-chlorophenyl)azetidine-1-carboxamide 858655-85-3P,

3-(4-Benzylpiperidin-1-yl)-N-(2-fluorophenyl)azetidine-1-carboxamide
858655-86-4P, 3-(4-Benzylpiperidin-1-yl)-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858655-87-5P,
3-(4-Benzylpiperidin-1-yl)-N-cyclopentylazetidine-1-carboxamide
858655-88-6P, 3-(4-Benzylpiperidin-1-yl)-N-(2,4-dimethylphenyl)azetidine-1-carboxamide 858655-89-7P,
3-(4-Benzylpiperidin-1-yl)-N-(3,5-dichlorophenyl)azetidine-1-carboxamide
858655-90-0P, 3-(4-Benzylpiperidin-1-yl)-N-[3,5-bis(trifluoromethyl)phenyl]azetidine-1-carboxamide 858655-91-1P,
3-(4-Benzylpiperidin-1-yl)-N-(3-phenoxyphenyl)azetidine-1-carboxamide
858655-92-2P, 3-(4-Benzylpiperidin-1-yl)-N-(3,5-difluorophenyl)azetidine-1-carboxamide 858655-93-3P,
3-(4-Benzylpiperidin-1-yl)-N-(4-methoxyphenyl)azetidine-1-carboxamide
858655-94-4P, 3-(4-Benzylpiperidin-1-yl)-N-(3,5-dimethylphenyl)azetidine-1-carboxamide 858655-95-5P,
3-(4-Benzylpiperidin-1-yl)-N-(3-fluorophenyl)azetidine-1-carboxamide
858655-96-6P, Methyl 3-([3-(4-benzylpiperidin-1-yl)azetidin-1-yl]carbonyl)amino)benzoate 858655-97-7P, 3-(4-Benzylpiperidin-1-yl)-N-[3-(methylsulfonyl)phenyl]azetidine-1-carboxamide
858655-98-8P, 3-(4-Benzylpiperidin-1-yl)-N-(2-methylphenyl)azetidine-1-carboxamide 858655-99-9P,
3-(2,3-Dihydro-1H-indol-1-yl)-N-ethylazetidine-1-carboxamide
858656-00-5P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-propylazetidine-1-carboxamide
858656-01-6P 858656-02-7P,
3-(2,3-Dihydro-1H-indol-1-yl)-N-hexylazetidine-1-carboxamide
858656-03-8P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(4-fluorophenyl)azetidine-1-carboxamide 858656-04-9P,
3-(2,3-Dihydro-1H-indol-1-yl)-N-(3-methylphenyl)azetidine-1-carboxamide
858656-05-0P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858656-06-1P,
3-(2,3-Dihydro-1H-indol-1-yl)-N-(4-isopropylphenyl)azetidine-1-carboxamide
858656-07-2P, N-(3-Chlorophenyl)-3-(2,3-dihydro-1H-indol-1-yl)azetidine-1-carboxamide 858656-08-3P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(2,5-dimethylphenyl)azetidine-1-carboxamide 858656-09-4P,
N-(4-Chlorophenyl)-3-(2,3-dihydro-1H-indol-1-yl)azetidine-1-carboxamide
858656-10-7P, N-Benzyl-3-(2,3-dihydro-1H-indol-1-yl)azetidine-1-carboxamide
858656-11-8P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(1-naphthyl)azetidine-1-carboxamide 858656-12-9P,
3-(2,3-Dihydro-1H-indol-1-yl)-N-(2-naphthyl)azetidine-1-carboxamide
858656-13-0P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858656-14-1P,
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858656-15-2P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(4-methylphenyl)azetidine-1-carboxamide 858656-16-3P,
N-Cyclohexyl-3-(2,3-dihydro-1H-indol-1-yl)azetidine-1-carboxamide
858656-17-4P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(2,6-dimethylphenyl)azetidine-1-carboxamide 858656-18-5P,
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858656-19-6P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(2-ethylphenyl)azetidine-1-carboxamide 858656-20-9P,
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858656-21-0P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(2-phenylethyl)azetidine-1-carboxamide 858656-22-1P,
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858656-23-2P, N-(2-Chlorophenyl)-3-(2,3-dihydro-1H-indol-1-yl)azetidine-1-carboxamide
858656-24-3P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(2-fluorophenyl)azetidine-1-carboxamide 858656-25-4P,
N-Cyclopentyl-3-(2,3-dihydro-1H-indol-1-yl)azetidine-1-carboxamide
858656-26-5P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(2,4-dimethylphenyl)azetidine-1-carboxamide 858656-27-6P,

N-(3,5-Difluorophenyl)-3-(2,3-dihydro-1H-indol-1-yl)azetidine-1-carboxamide 858656-28-7P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(4-methoxyphenyl)azetidine-1-carboxamide 858656-29-8P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(3,5-dimethylphenyl)azetidine-1-carboxamide 858656-30-1P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(3-fluorophenyl)azetidine-1-carboxamide 858656-31-2P, Methyl 3-([3-(2,3-dihydro-1H-indol-1-yl)azetidin-1-yl]carbonyl)amino)benzoate 858656-32-3P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-[3-(methylsulfanyl)phenyl]azetidine-1-carboxamide 858656-33-4P, 3-(2,3-Dihydro-1H-indol-1-yl)-N-(2-methylphenyl)azetidine-1-carboxamide 858656-34-5P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-propylazetidine-1-carboxamide 858656-35-6P 858656-36-7P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-hexylazetidine-1-carboxamide 858656-37-8P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(4-fluorophenyl)azetidine-1-carboxamide 858656-38-9P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(3-methylphenyl)azetidine-1-carboxamide 858656-39-0P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(2-(thien-2-yl)ethyl)azetidine-1-carboxamide 858656-40-3P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(4-isopropylphenyl)azetidine-1-carboxamide 858656-41-4P, N-(3-Chlorophenyl)-3-(3,4-dihydroquinolin-1(2H)-yl)azetidine-1-carboxamide 858656-42-5P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(2,5-dimethylphenyl)azetidine-1-carboxamide 858656-43-6P, N-(4-Chlorophenyl)-3-(3,4-dihydroquinolin-1(2H)-yl)azetidine-1-carboxamide 858656-44-7P, N-Benzyl-3-(3,4-dihydroquinolin-1(2H)-yl)azetidine-1-carboxamide 858656-45-8P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(1-naphthyl)azetidine-1-carboxamide 858656-46-9P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(2-naphthyl)azetidine-1-carboxamide 858656-47-0P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-[1-(1-naphthyl)ethyl]azetidine-1-carboxamide 858656-48-1P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(3,4-dimethylphenyl)azetidine-1-carboxamide 858656-49-2P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(4-methylphenyl)azetidine-1-carboxamide 858656-50-5P, N-Cyclohexyl-3-(3,4-dihydroquinolin-1(2H)-yl)azetidine-1-carboxamide 858656-51-6P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(2,6-dimethylphenyl)azetidine-1-carboxamide 858656-52-7P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(2-ethoxyphenyl)azetidine-1-carboxamide 858656-53-8P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(2-ethylphenyl)azetidine-1-carboxamide 858656-54-9P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(4-ethoxyphenyl)azetidine-1-carboxamide 858656-56-1P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(2-phenylethyl)azetidine-1-carboxamide 858656-57-2P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-phenylazetidine-1-carboxamide 858656-59-4P, N-(2-Chlorophenyl)-3-(3,4-dihydroquinolin-1(2H)-yl)azetidine-1-carboxamide 858656-61-8P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(2-fluorophenyl)azetidine-1-carboxamide 858656-64-1P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-[3-(trifluoromethyl)phenyl]azetidine-1-carboxamide 858656-65-2P, N-Cyclopentyl-3-(3,4-dihydroquinolin-1(2H)-yl)azetidine-1-carboxamide 858656-66-3P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(2,4-dimethylphenyl)azetidine-1-carboxamide 858656-67-4P, N-(3,5-Dichlorophenyl)-3-(3,4-dihydroquinolin-1(2H)-yl)azetidine-1-carboxamide 858656-68-5P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-(3,4-dihydroquinolin-1(2H)-yl)azetidine-1-carboxamide 858656-69-6P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(3-phenoxyphenyl)azetidine-1-carboxamide 858656-70-9P, N-(3,5-Difluorophenyl)-3-(3,4-dihydroquinolin-1(2H)-yl)azetidine-1-carboxamide 858656-71-0P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(4-methoxyphenyl)azetidine-1-carboxamide 858656-72-1P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(3,5-dimethylphenyl)azetidine-1-carboxamide 858656-73-2P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(3-

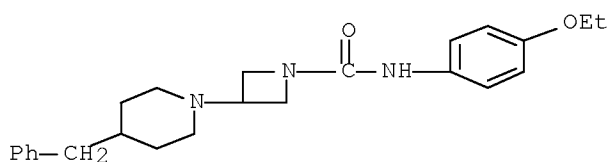
fluorophenyl)azetidine-1-carboxamide 858656-74-3P, Methyl
3-([3-(3,4-dihydroquinolin-1(2H)-yl)azetidin-1-yl]carbonyl)amino)benzoate
858656-75-4P, 3-(3,4-Dihydroquinolin-1(2H)-yl)-N-[3-
(methylsulfanyl)phenyl]azetidine-1-carboxamide 858656-76-5P,
3-(3,4-Dihydroquinolin-1(2H)-yl)-N-(2-methylphenyl)azetidine-1-carboxamide
858656-78-7P, 3-(4-Benzylpiperazin-1-yl)-N-[3,5-
bis(trifluoromethyl)phenyl]azetidine-1-carboxamide 858656-80-1P
858656-81-2P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-[(1-methyl-
1H-pyrrol-2-yl)methyl]piperazin-1-yl]azetidine-1-carboxamide
858656-82-3P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-[(1,5-
dimethyl-3-oxo-2-phenyl-2,3-dihydro-1H-pyrazol-4-yl)methyl]piperazin-1-
yl]azetidine-1-carboxamide 858656-83-4P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-[(5-methyl-2-furyl)methyl]piperazin-1-
yl]azetidine-1-carboxamide 858656-84-5P, [5-([4-[1-([3,5-
Bis(trifluoromethyl)phenyl)amino]carbonyl)azetidin-3-yl]piperazin-1-
yl)methyl]-2-furyl]methyl acetate 858656-85-6P,
N-[3,5-Bis(trifluoromethyl)phenyl]-3-(4-[[5-(hydroxymethyl)-2-
furyl]methyl]piperazin-1-yl)azetidine-1-carboxamide 858656-86-7P
, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(2-methoxybenzyl)piperazin-1-
yl]azetidine-1-carboxamide 858656-87-8P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-(2,3-dimethoxybenzyl)piperazin-1-
yl]azetidine-1-carboxamide 858656-88-9P,
N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(2,4-dimethoxybenzyl)piperazin-1-
yl]azetidine-1-carboxamide 858656-89-0P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-(2,4,6-trimethoxybenzyl)piperazin-1-
yl]azetidine-1-carboxamide 858656-90-3P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-(2,5-dimethoxybenzyl)piperazin-1-
yl]azetidine-1-carboxamide 858656-91-4P, [2-([4-[1-([3,5-
Bis(trifluoromethyl)phenyl)amino]carbonyl)azetidin-3-yl]piperazin-1-
yl)methyl]phenoxy]acetic acid 858656-92-5P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-(2-(trifluoromethyl)benzyl)piperazin-1-
yl]azetidine-1-carboxamide 858656-93-6P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-(2-methylbenzyl)piperazin-1-yl]azetidine-
1-carboxamide 858656-94-7P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-
[4-(3-cyanobenzyl)piperazin-1-yl]azetidine-1-carboxamide
858656-95-8P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(3-
fluorobenzyl)piperazin-1-yl]azetidine-1-carboxamide 858656-96-9P
, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(3-fluoro-4-
methoxybenzyl)piperazin-1-yl]azetidine-1-carboxamide 858656-97-0P
, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(3-phenoxybenzyl)piperazin-1-
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Bis(trifluoromethyl)phenyl]-3-[4-(3-methoxybenzyl)piperazin-1-yl]azetidine-
1-carboxamide 858656-99-2P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-
[4-(3,4-dimethoxybenzyl)piperazin-1-yl]azetidine-1-carboxamide
858657-00-8P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(3,4,5-
trimethoxybenzyl)piperazin-1-yl]azetidine-1-carboxamide
858657-01-9P, 3-[4-[4-(Benzyloxy)-3-methoxybenzyl]piperazin-1-yl]-
N-[3,5-bis(trifluoromethyl)phenyl]azetidine-1-carboxamide
858657-02-0P, 3-[4-[3-(Benzyloxy)benzyl]piperazin-1-yl]-N-[3,5-
bis(trifluoromethyl)phenyl]azetidine-1-carboxamide 858657-03-1P,
N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(3-hydroxybenzyl)piperazin-1-
yl]azetidine-1-carboxamide 858657-04-2P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-(3-hydroxy-4-methoxybenzyl)piperazin-1-
yl]azetidine-1-carboxamide 858657-05-3P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-[3-(trifluoromethyl)benzyl]piperazin-1-
yl]azetidine-1-carboxamide 858657-06-4P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-(3-methylbenzyl)piperazin-1-yl]azetidine-
1-carboxamide 858657-07-5P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-
[4-(4-cyanobenzyl)piperazin-1-yl]azetidine-1-carboxamide
858657-08-6P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(4-

fluorobenzyl)piperazin-1-yl]azetidine-1-carboxamide 858657-09-7P,
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bis(trifluoromethyl)phenyl]azetidine-1-carboxamide 858657-10-0P,
N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-[4-(dimethylamino)benzyl]piperazin-
1-yl]azetidine-1-carboxamide 858657-11-1P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-[4-(diethylamino)benzyl]piperazin-1-
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Bis(trifluoromethyl)phenyl]-3-[4-(4-phenoxybenzyl)piperazin-1-yl]azetidine-
1-carboxamide 858657-13-3P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-
[4-(4-methoxybenzyl)piperazin-1-yl]azetidine-1-carboxamide
858657-14-4P, 3-[4-[4-(Benzyloxy)benzyl]piperazin-1-yl]-N-[3,5-
bis(trifluoromethyl)phenyl]azetidine-1-carboxamide 858657-15-5P,
N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(1H-imidazol-2-ylmethyl)piperazin-
1-yl]azetidine-1-carboxamide 858657-16-6P, N-[3,5-
Bis(trifluoromethyl)phenyl]-3-[4-(1-naphthylmethyl)piperazin-1-
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yl]azetidine-1-carboxamide 858657-18-8P, 3-[4-[3,4-
Bis(benzyloxy)benzyl]piperazin-1-yl]-N-[3,5-bis(trifluoromethyl)phenyl]aze
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yl]azetidine-1-carboxamide 858657-20-2P, N-[3,5-
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Bis(trifluoromethyl)phenyl]-3-[4-[(3-methylthien-2-yl)methyl]piperazin-1-
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Bis(trifluoromethyl)phenyl]-3-[4-[(4-bromothien-2-yl)methyl]piperazin-1-
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1-carboxamide 858657-27-9P 858657-28-0P, Methyl
4-([4-[1-([3,5-bis(trifluoromethyl)phenyl]amino)carbonyl]azetidin-3-
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N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-[4-(trifluoromethyl)benzyl]piperaz
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Bis(trifluoromethyl)phenyl]amino)carbonyl]azetidin-3-yl]piperazin-1-
yl]acetic acid 858657-33-7P, N-[3,5-Bis(trifluoromethyl)phenyl]-
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858657-34-8P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-(4-
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(4-propylpiperazin-1-yl)azetidine-1-carboxamide 858657-39-3P,
N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-[3-(methylsulfanyl)propyl]piperazi
n-1-yl]azetidine-1-carboxamide 858657-40-6P,
N-[3,5-Bis(trifluoromethyl)phenyl]-3-(4-butylpiperazin-1-yl)azetidine-1-
carboxamide 858657-41-7P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-

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 858657-42-8P, N-[3,5-Bis(trifluoromethyl)phenyl]-3-[4-(3-nitrobenzyl)piperazin-1-yl]azetidine-1-carboxamide 858657-43-9P,
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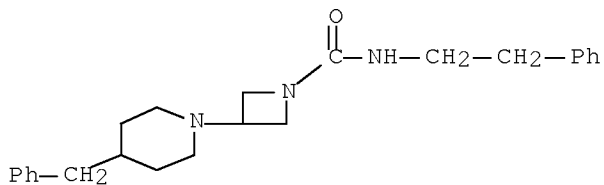
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 , Methyl 2-([4-[1-([3,5-bis(trifluoromethyl)phenyl]amino)carbonyl]azetidi
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 RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
 (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
 (Uses)

(preparation of azetidine ring compds. as EDG-5 antagonists for treatment
 of
 hypertension, myocardial infarction, etc.)
 RN 858655-81-9 CAPLUS
 CN 1-Azetidinecarboxamide, N-(4-ethoxyphenyl)-3-[4-(phenylmethyl)-1-
 piperidinyl]- (CA INDEX NAME)



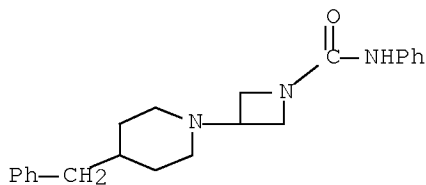
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CN 1-Azetidinecarboxamide, N-(2-phenylethyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



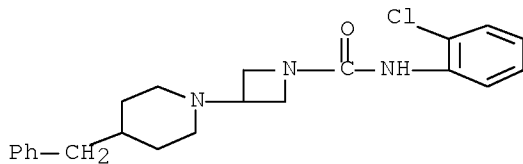
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CN 1-Azetidinecarboxamide, N-phenyl-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



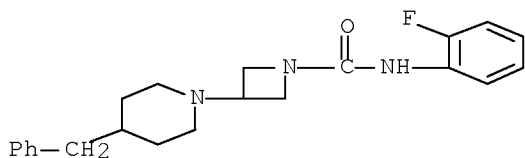
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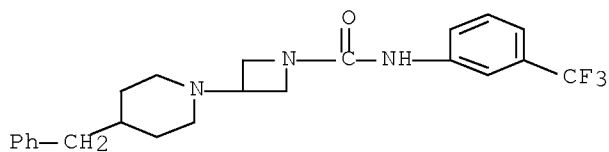
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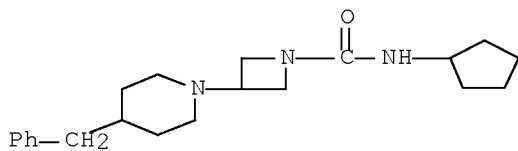
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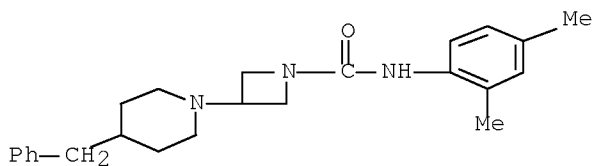
RN 858655-87-5 CAPLUS

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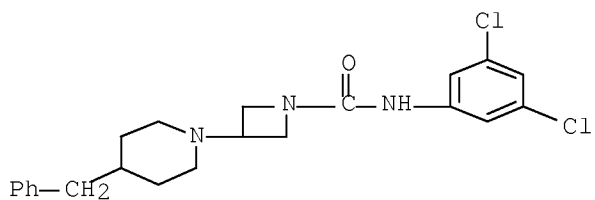
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CN 1-Azetidinecarboxamide, N-(2,4-dimethylphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



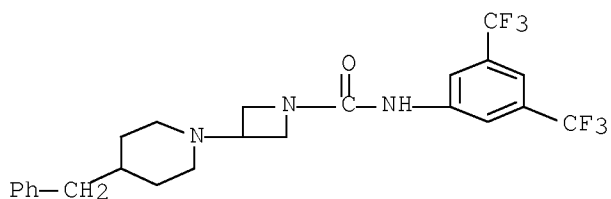
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CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



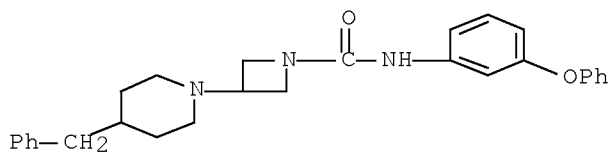
RN 858655-90-0 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



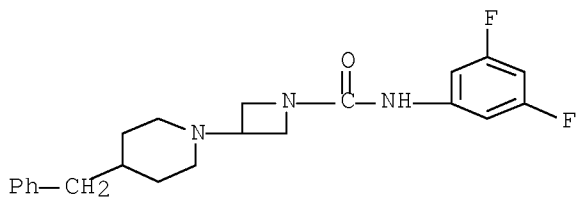
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CN 1-Azetidinecarboxamide, N-(3-phenoxyphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



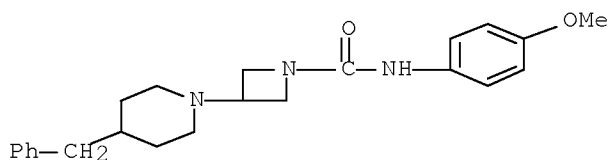
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CN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



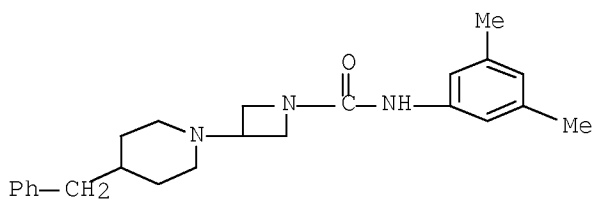
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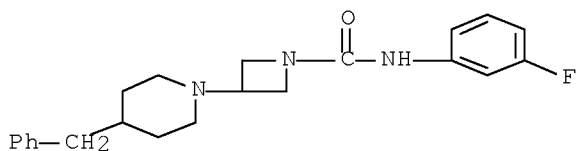
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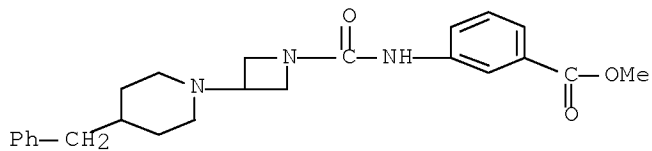
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CN 1-Azetidinecarboxamide, N-(3-fluorophenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



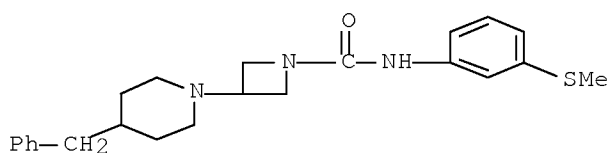
RN 858655-96-6 CAPLUS

CN Benzoic acid, 3-[[[3-[4-(phenylmethyl)-1-piperidinyl]-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



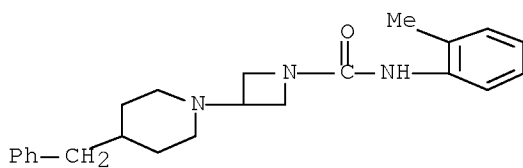
RN 858655-97-7 CAPLUS

CN 1-Azetidinecarboxamide, N-[3-(methylthio)phenyl]-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



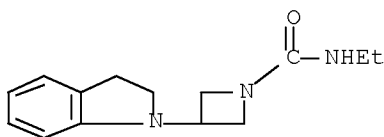
RN 858655-98-8 CAPLUS

CN 1-Azetidinecarboxamide, N-(2-methylphenyl)-3-[4-(phenylmethyl)-1-piperidinyl]- (CA INDEX NAME)



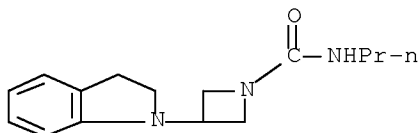
RN 858655-99-9 CAPLUS

CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-ethyl- (CA INDEX NAME)



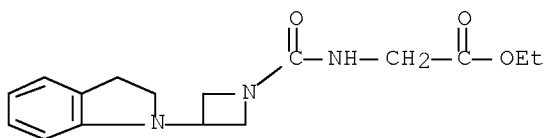
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-propyl- (CA INDEX NAME)



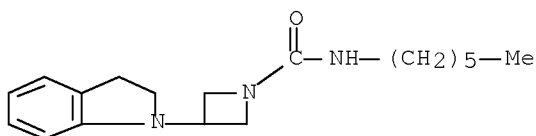
RN 858656-01-6 CAPLUS

CN Glycine, N-[[3-(2,3-dihydro-1H-indol-1-yl)-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)



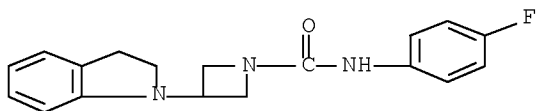
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-hexyl- (CA INDEX NAME)



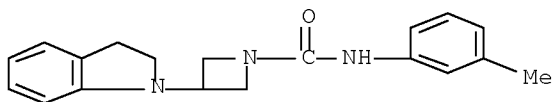
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(4-fluorophenyl)- (CA INDEX NAME)



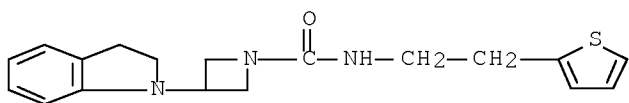
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(3-methylphenyl)- (CA INDEX NAME)



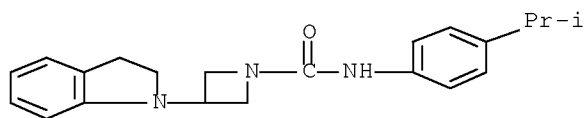
RN 858656-05-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-[2-(2-thienyl)ethyl]- (CA INDEX NAME)



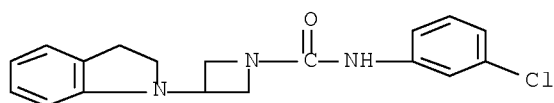
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-[4-(1-methylethyl)phenyl]- (CA INDEX NAME)



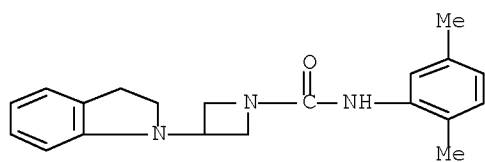
RN 858656-07-2 CAPLUS

CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-(2,3-dihydro-1H-indol-1-yl)- (CA INDEX NAME)



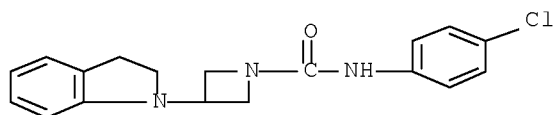
RN 858656-08-3 CAPLUS

CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(2,5-dimethylphenyl)- (CA INDEX NAME)



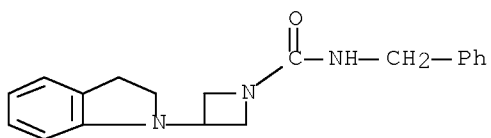
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CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-(2,3-dihydro-1H-indol-1-yl)- (CA INDEX NAME)



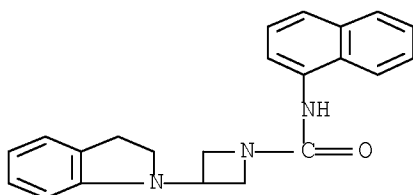
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(phenylmethyl)- (CA INDEX NAME)



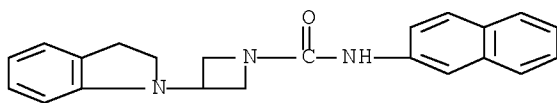
RN 858656-11-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-1-naphthalenyl-
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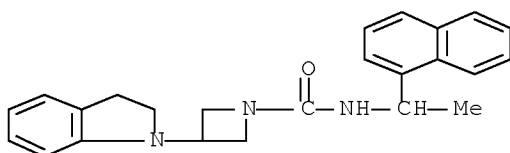
RN 858656-12-9 CAPLUS

CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-2-naphthalenyl-
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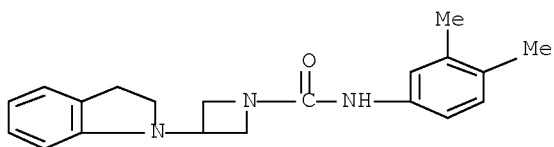
RN 858656-13-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-[1-(1-naphthalenyl)ethyl]-
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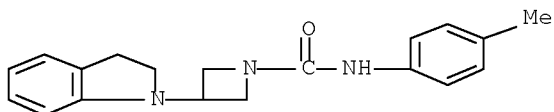
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(3,4-dimethylphenyl)-
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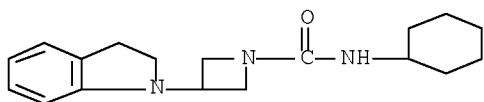
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(4-methylphenyl)-
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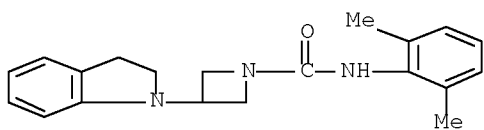
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CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(2,3-dihydro-1H-indol-1-yl)- (CA
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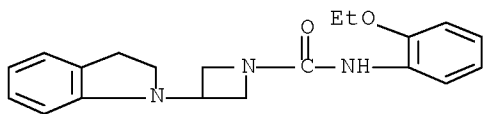
RN 858656-17-4 CAPLUS

CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(2,6-
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RN 858656-18-5 CAPLUS

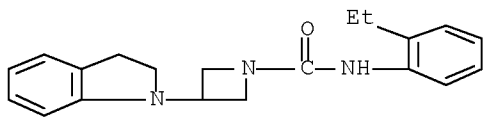
CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(2-ethoxyphenyl)-
(CA INDEX NAME)



RN 858656-19-6 CAPLUS

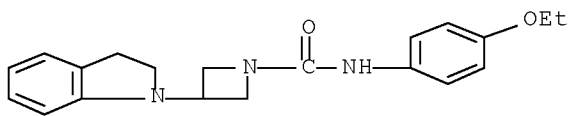
CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(2-ethylphenyl)-

(CA INDEX NAME)



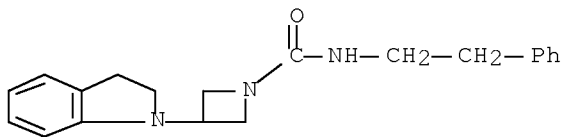
RN 858656-20-9 CAPLUS

CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(4-ethoxyphenyl)-
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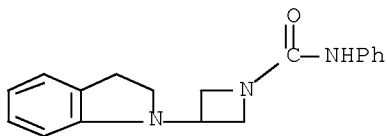
RN 858656-21-0 CAPLUS

CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(2-phenylethyl)-
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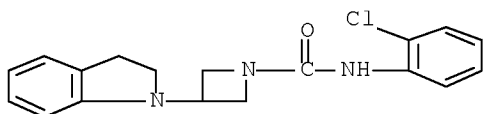
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-phenyl- (CA INDEX
NAME)



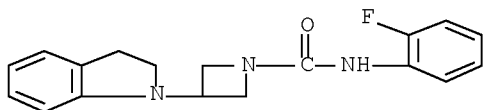
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CN 1-Azetidinecarboxamide, N-(2-chlorophenyl)-3-(2,3-dihydro-1H-indol-1-yl)-
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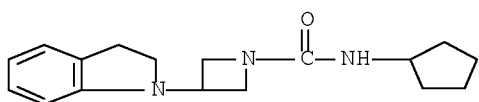
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(2-fluorophenyl)-
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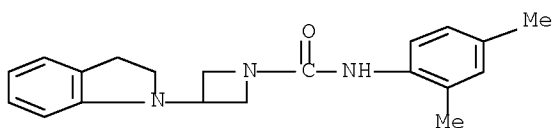
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CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(2,3-dihydro-1H-indol-1-yl)- (CA
INDEX NAME)



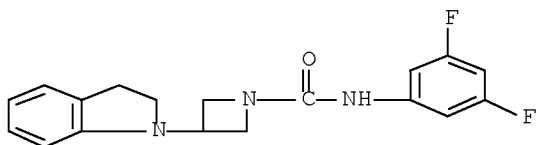
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(2,4-
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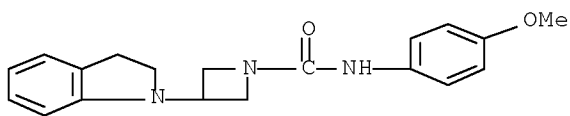
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CN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-(2,3-dihydro-1H-indol-1-
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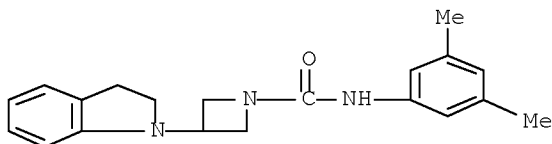
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(4-methoxyphenyl)-
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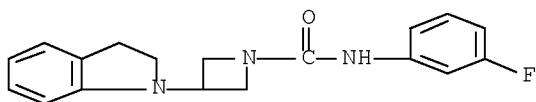
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(3,5-dimethylphenyl)- (CA INDEX NAME)



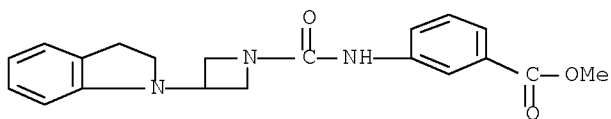
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CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(3-fluorophenyl)- (CA INDEX NAME)



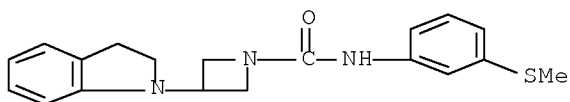
RN 858656-31-2 CAPLUS

CN Benzoic acid, 3-[[[3-(2,3-dihydro-1H-indol-1-yl)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



RN 858656-32-3 CAPLUS

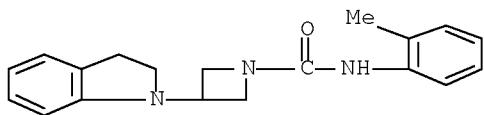
CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-[3-(methylthio)phenyl]- (CA INDEX NAME)



RN 858656-33-4 CAPLUS

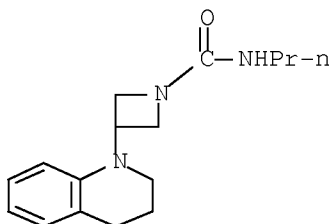
CN 1-Azetidinecarboxamide, 3-(2,3-dihydro-1H-indol-1-yl)-N-(2-methylphenyl)-

(CA INDEX NAME)



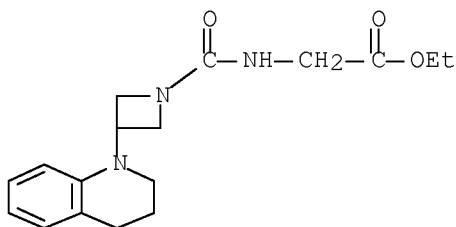
RN 858656-34-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-propyl- (CA INDEX NAME)



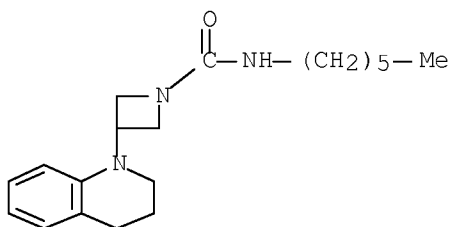
RN 858656-35-6 CAPLUS

CN Glycine, N-[[3-(3,4-dihydro-1(2H)-quinolinyl)-1-azetidiny]carbonyl]-, ethyl ester (CA INDEX NAME)



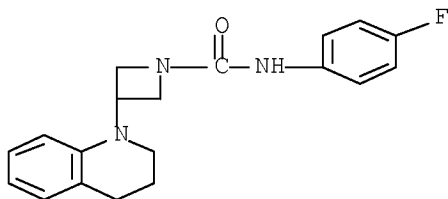
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-hexyl- (CA INDEX NAME)



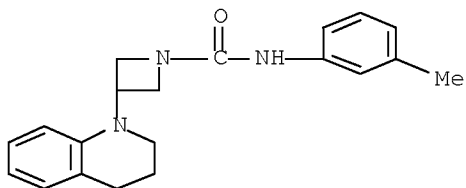
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(4-fluorophenyl)- (CA INDEX NAME)



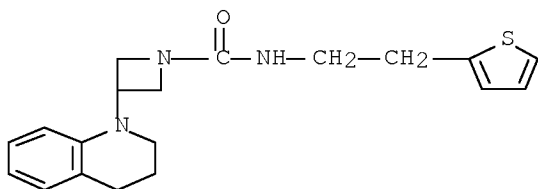
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(3-methylphenyl)- (CA INDEX NAME)



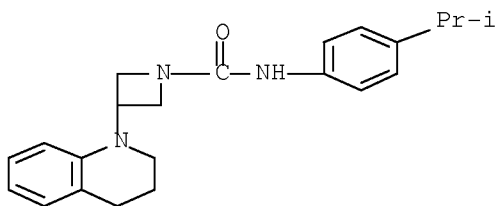
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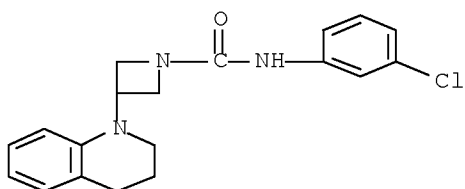


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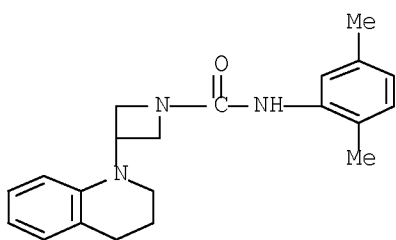
CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-[4-(1-methylethyl)phenyl]- (CA INDEX NAME)



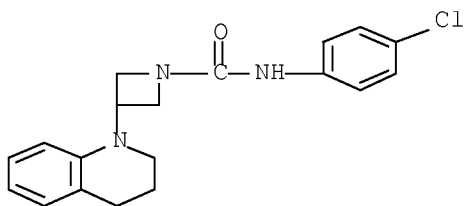
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CN 1-Azetidinecarboxamide, N-(3-chlorophenyl)-3-(3,4-dihydro-1(2H)-quinolinyl)- (CA INDEX NAME)



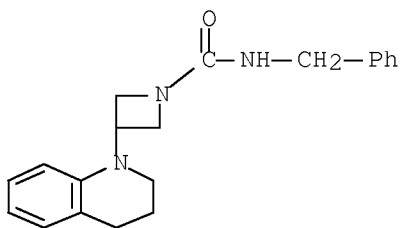
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(2,5-dimethylphenyl)- (CA INDEX NAME)



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CN 1-Azetidinecarboxamide, N-(4-chlorophenyl)-3-(3,4-dihydro-1(2H)-quinolinyl)- (CA INDEX NAME)

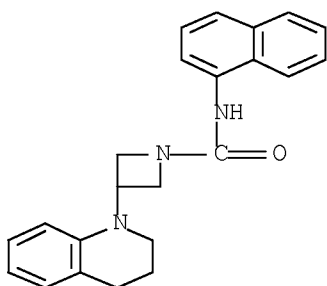


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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(phenylmethyl)- (CA INDEX NAME)



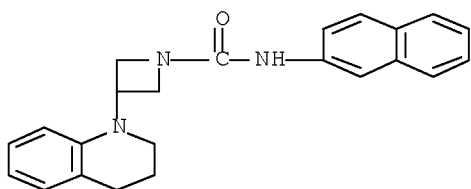
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-1-naphthalenyl-
(CA INDEX NAME)



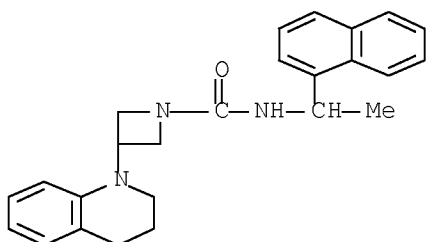
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-2-naphthalenyl-
(CA INDEX NAME)



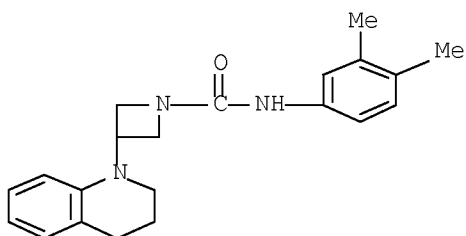
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-[1-(1-naphthalenyl)ethyl]- (CA INDEX NAME)



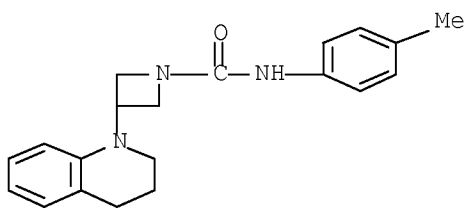
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(3,4-dimethylphenyl)- (CA INDEX NAME)



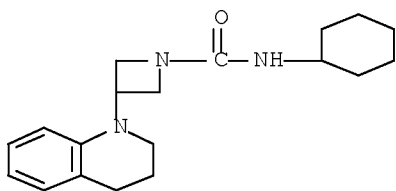
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(4-methylphenyl)- (CA INDEX NAME)



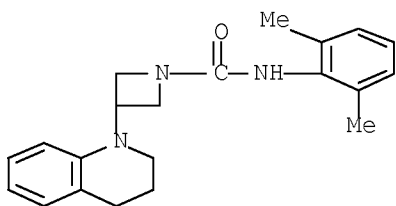
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CN 1-Azetidinecarboxamide, N-cyclohexyl-3-(3,4-dihydro-1(2H)-quinolinyl)- (CA INDEX NAME)



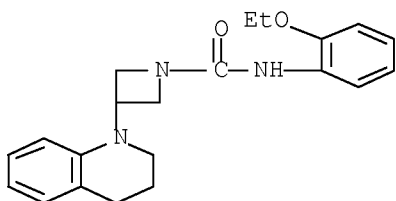
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(2,6-dimethylphenyl)- (CA INDEX NAME)



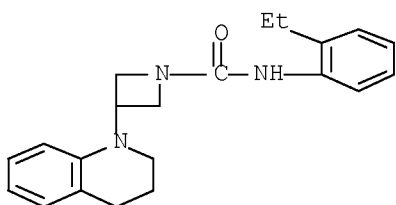
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(2-ethoxyphenyl)- (CA INDEX NAME)



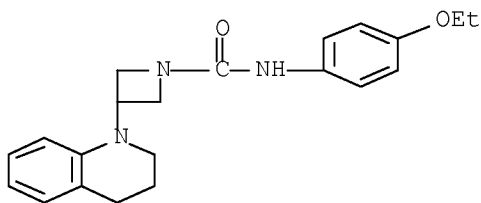
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(2-ethylphenyl)- (CA INDEX NAME)



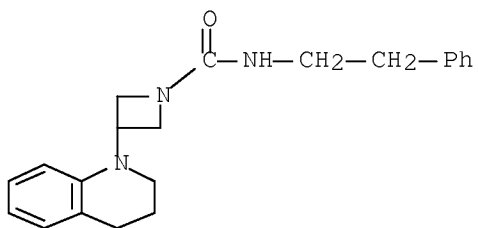
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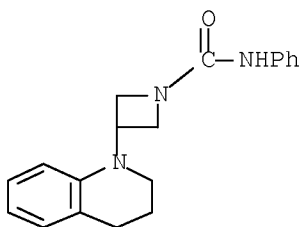


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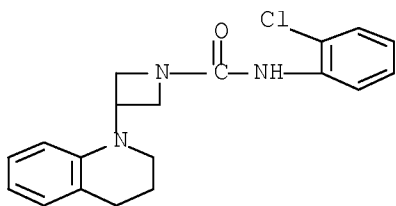
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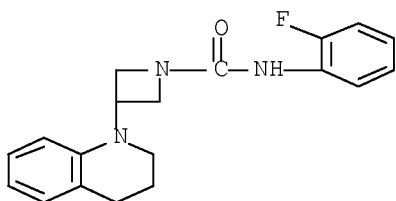
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 CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-phenyl- (CA INDEX NAME)



RN 858656-59-4 CAPLUS
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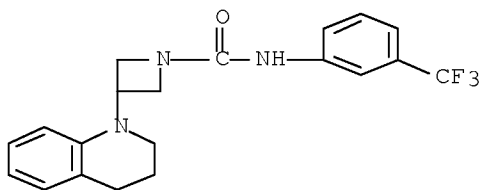


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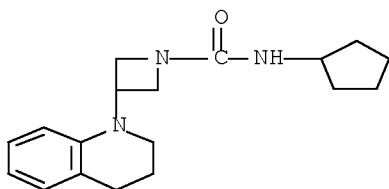
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



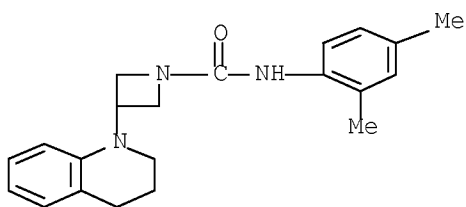
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CN 1-Azetidinecarboxamide, N-cyclopentyl-3-(3,4-dihydro-1(2H)-quinolinyl)- (CA INDEX NAME)



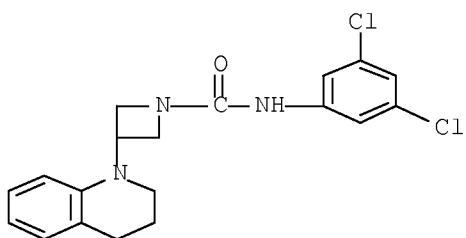
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(2,4-dimethylphenyl)- (CA INDEX NAME)

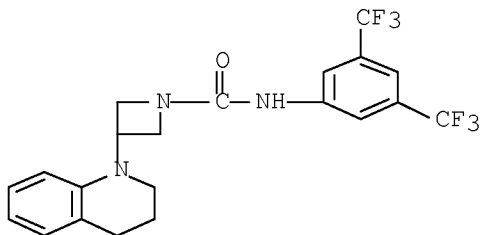


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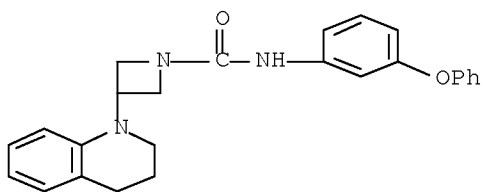
CN 1-Azetidinecarboxamide, N-(3,5-dichlorophenyl)-3-(3,4-dihydro-1(2H)-quinolinyl)- (CA INDEX NAME)



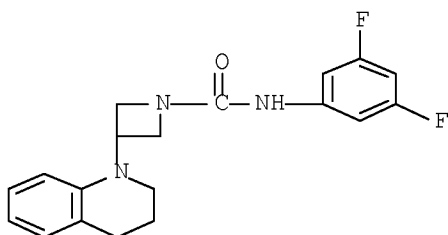
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(3,4-dihydro-1(2H)-quinolinyl)- (CA INDEX NAME)



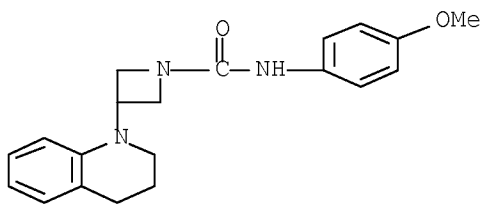
RN 858656-69-6 CAPLUS
CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(3-phenoxyphenyl)- (CA INDEX NAME)



RN 858656-70-9 CAPLUS
CN 1-Azetidinecarboxamide, N-(3,5-difluorophenyl)-3-(3,4-dihydro-1(2H)-quinolinyl)- (CA INDEX NAME)

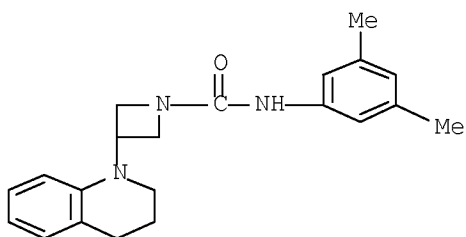


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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(4-methoxyphenyl)- (CA INDEX NAME)



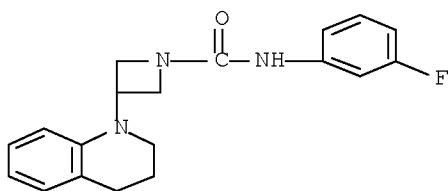
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(3,5-dimethylphenyl)- (CA INDEX NAME)



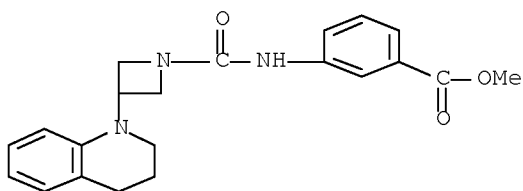
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(3-fluorophenyl)- (CA INDEX NAME)



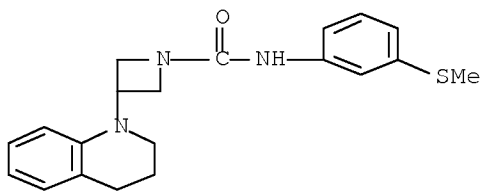
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CN Benzoic acid, 3-[[[3-(3,4-dihydro-1(2H)-quinolinyl)-1-azetidiny]carbonyl]amino]-, methyl ester (CA INDEX NAME)



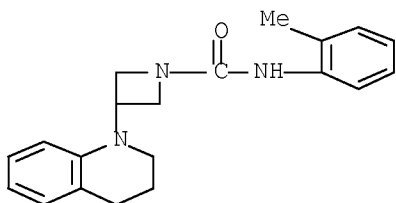
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CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-[3-(methylthio)phenyl]- (CA INDEX NAME)



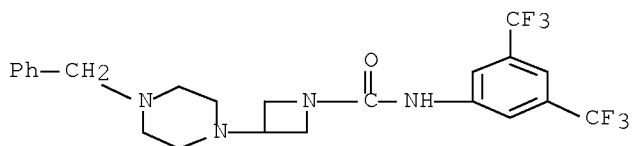
RN 858656-76-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-(3,4-dihydro-1(2H)-quinolinyl)-N-(2-methylphenyl)- (CA INDEX NAME)



RN 858656-78-7 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(phenylmethyl)-1-piperazinyl]- (CA INDEX NAME)



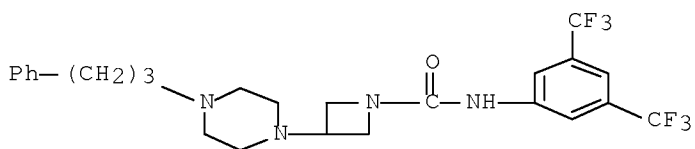
RN 858656-80-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(3-phenylpropyl)-1-piperazinyl]-, monoacetate (9CI) (CA INDEX NAME)

CM 1

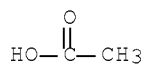
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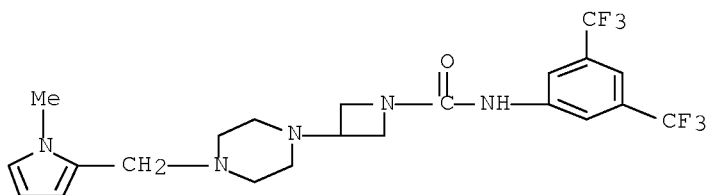


CM 2

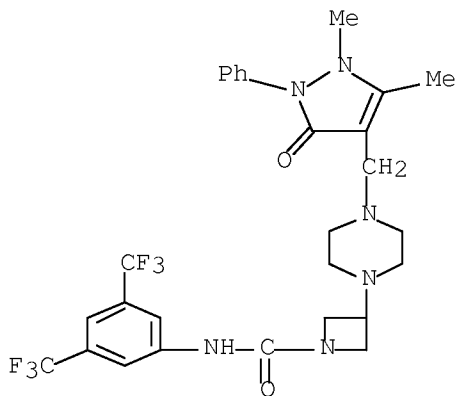
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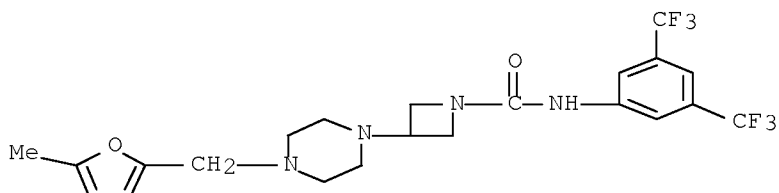
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(1-methyl-1H-pyrrol-2-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



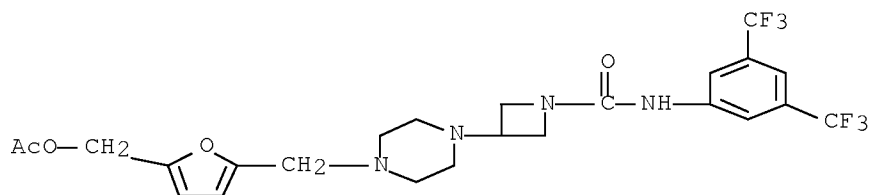
RN 858656-82-3 CAPLUS
CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(2,3-dihydro-1,5-dimethyl-3-oxo-2-phenyl-1H-pyrazol-4-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



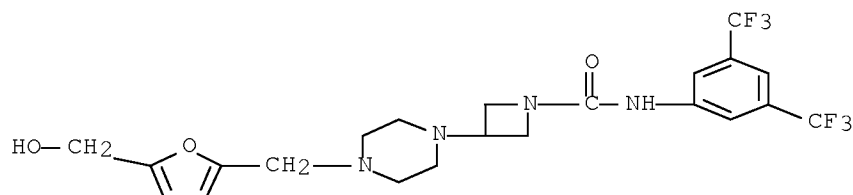
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(5-methyl-2-furanyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



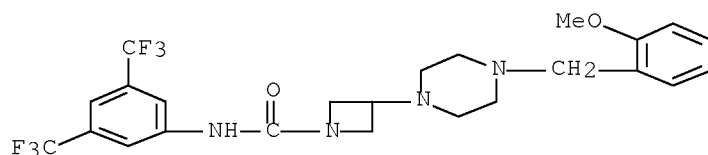
RN 858656-84-5 CAPLUS
 CN 1-Azetidinecarboxamide, 3-[4-[[5-[(acetyloxy)methyl]-2-furanyl]methyl]-1-piperazinyl]-N-[3,5-bis(trifluoromethyl)phenyl]- (CA INDEX NAME)



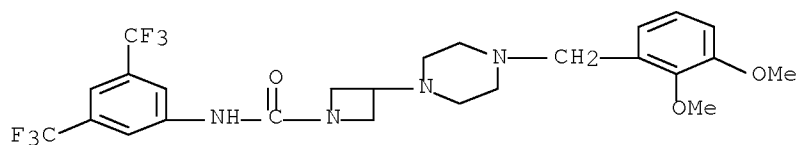
RN 858656-85-6 CAPLUS
 CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[5-(hydroxymethyl)-2-furanyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



RN 858656-86-7 CAPLUS
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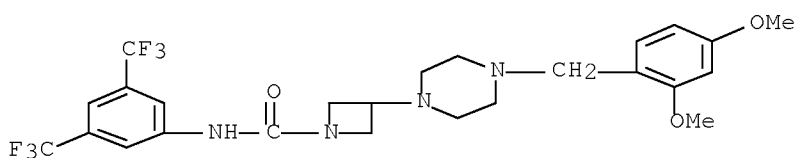


RN 858656-87-8 CAPLUS
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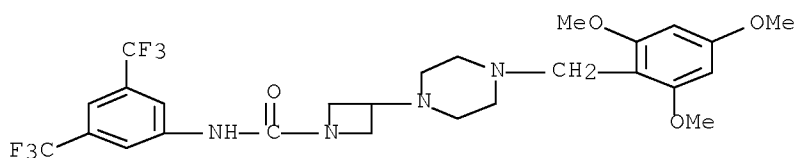
RN 858656-88-9 CAPLUS
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dimethoxyphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



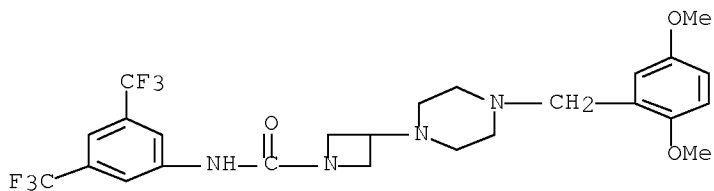
RN 858656-89-0 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(2,4,6-trimethoxyphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



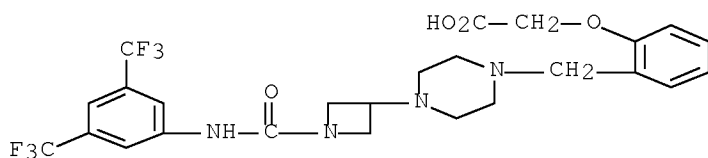
RN 858656-90-3 CAPLUS

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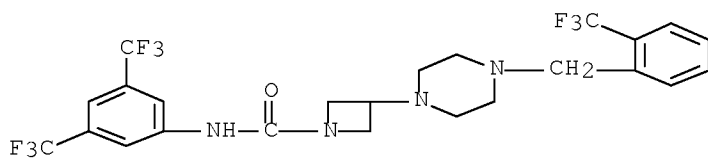
RN 858656-91-4 CAPLUS

CN Acetic acid, [2-[[4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]-1-piperazinyl]methyl]phenoxy]- (9CI) (CA INDEX NAME)



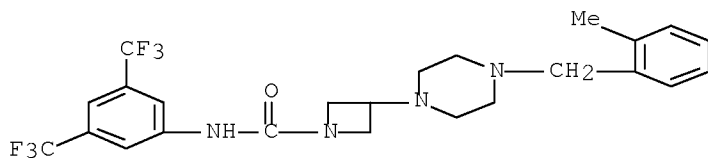
RN 858656-92-5 CAPLUS

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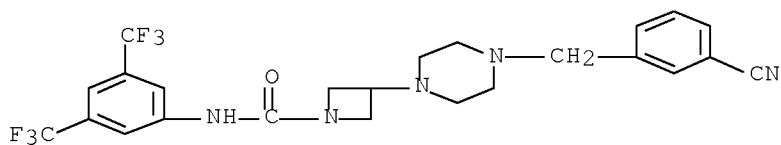
RN 858656-93-6 CAPLUS

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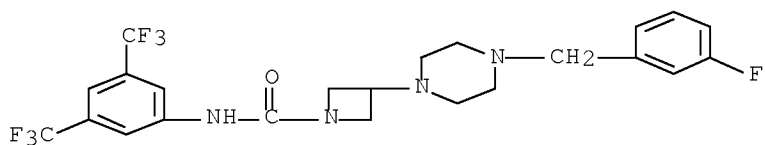
RN 858656-94-7 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-cyanophenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



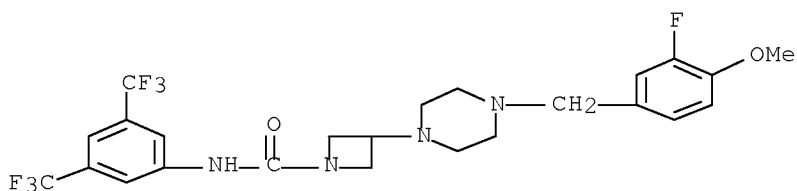
RN 858656-95-8 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-fluorophenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



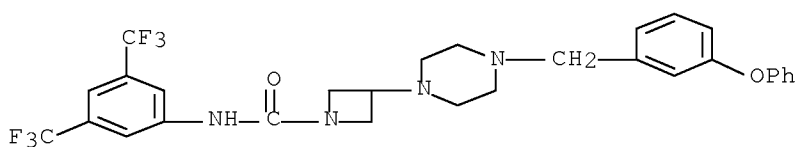
RN 858656-96-9 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-fluoro-4-methoxyphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



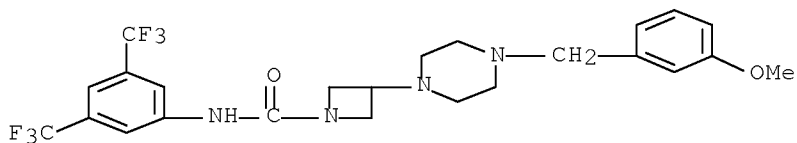
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-phenoxyphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



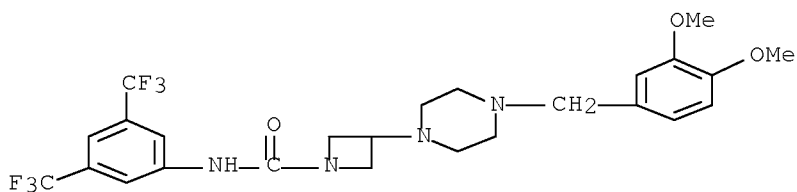
RN 858656-98-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-methoxyphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



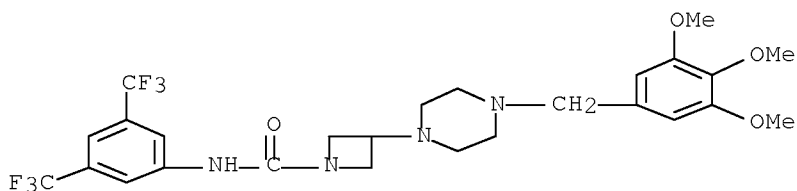
RN 858656-99-2 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3,4-dimethoxyphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



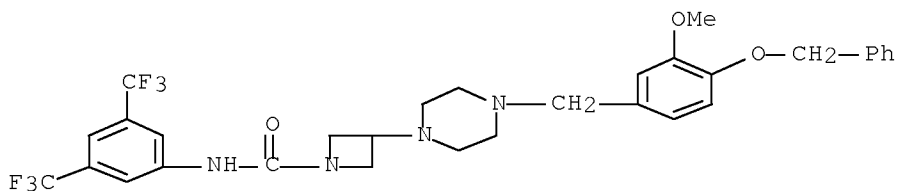
RN 858657-00-8 CAPLUS

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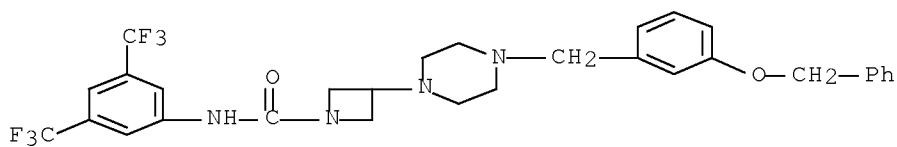
RN 858657-01-9 CAPLUS

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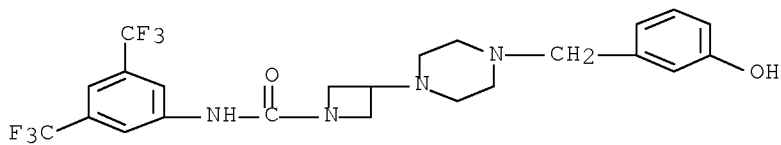
RN 858657-02-0 CAPLUS

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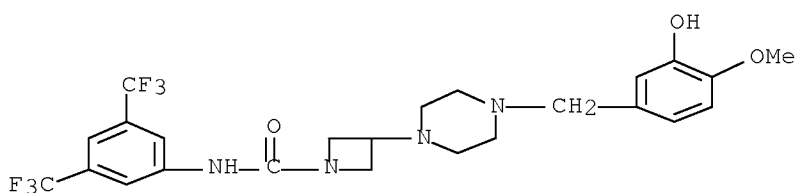
RN 858657-03-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[3-hydroxyphenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



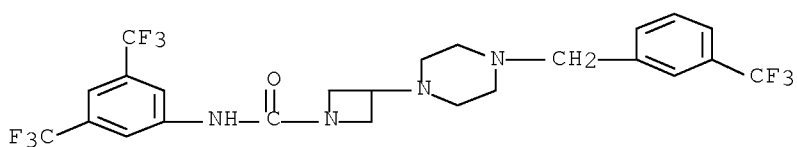
RN 858657-04-2 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[3-hydroxy-4-methoxyphenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



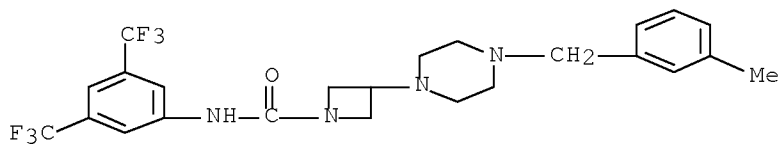
RN 858657-05-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-(trifluoromethyl)phenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



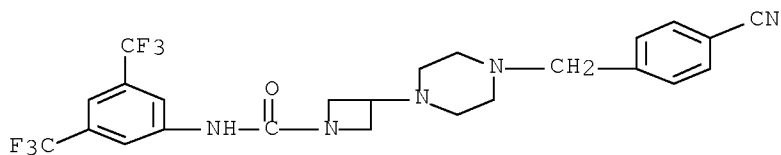
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-methylphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



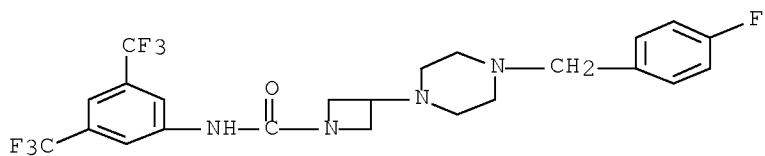
RN 858657-07-5 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(4-cyanophenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



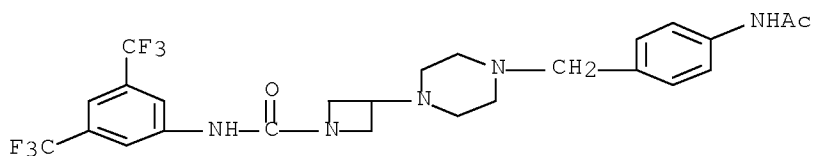
RN 858657-08-6 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(4-fluorophenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



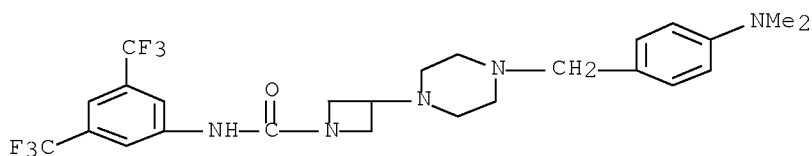
RN 858657-09-7 CAPLUS

CN 1-Azetidinecarboxamide, 3-[4-[[4-(acetylamino)phenyl]methyl]-1-piperazinyl]-N-[3,5-bis(trifluoromethyl)phenyl]- (CA INDEX NAME)



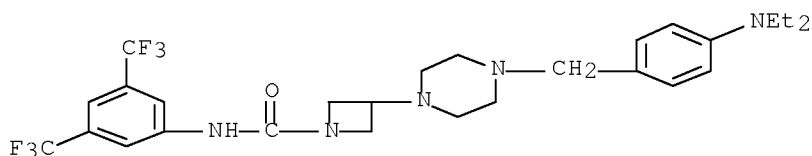
RN 858657-10-0 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(dimethylamino)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



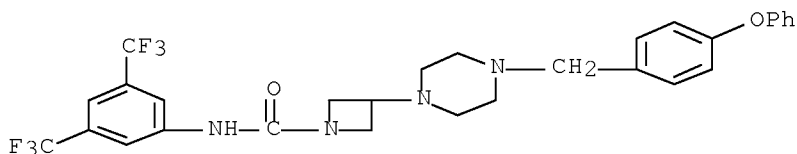
RN 858657-11-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(diethylamino)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



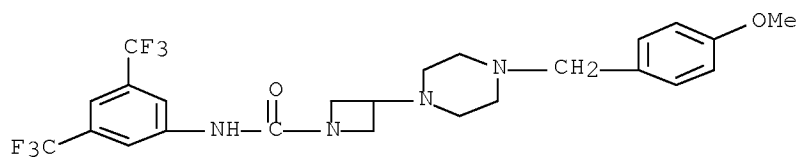
RN 858657-12-2 CAPLUS

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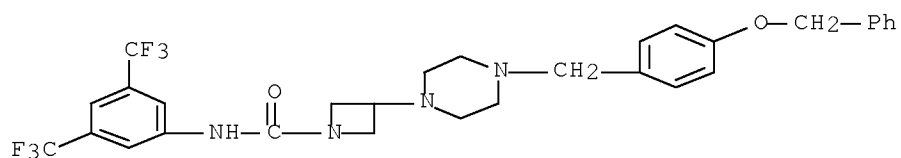
RN 858657-13-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(4-methoxyphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



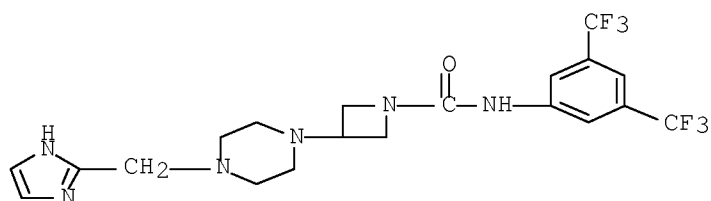
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(phenylmethoxy)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



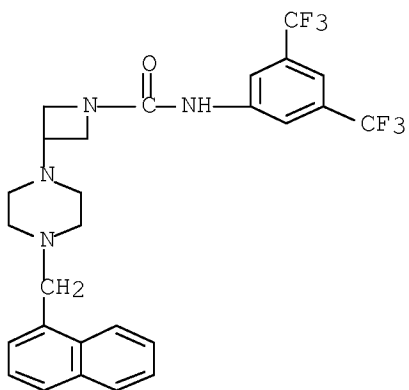
RN 858657-15-5 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(1H-imidazol-2-ylmethyl)-1-piperazinyl]- (CA INDEX NAME)



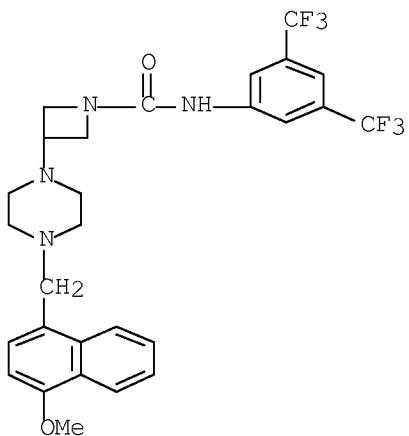
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(1-naphthalenylmethyl)-1-piperazinyl]- (CA INDEX NAME)



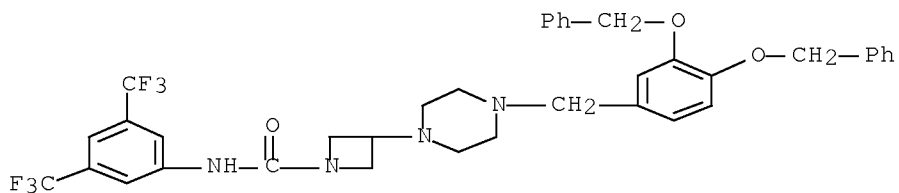
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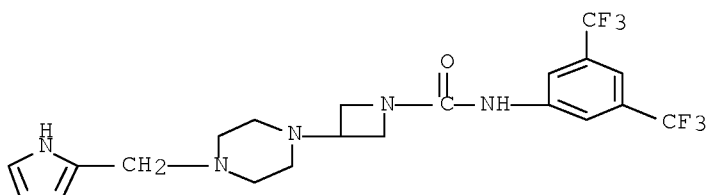
RN 858657-18-8 CAPLUS

CN 1-Azetidinecarboxamide, 3-[4-[[3,4-bis(phenylmethoxy)phenyl]methyl]-1-piperazinyl]-N-[3,5-bis(trifluoromethyl)phenyl]- (CA INDEX NAME)



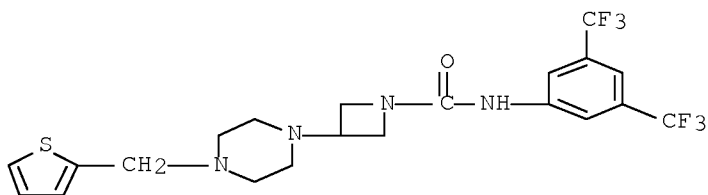
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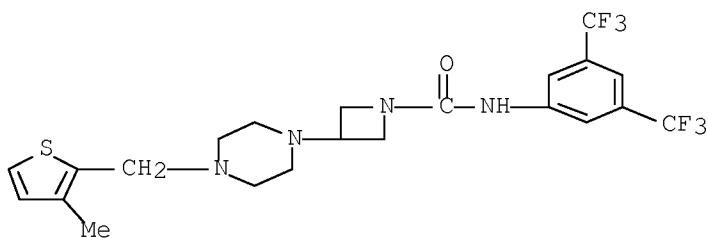
RN 858657-20-2 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(2-thienylmethyl)-1-piperazinyl]- (CA INDEX NAME)



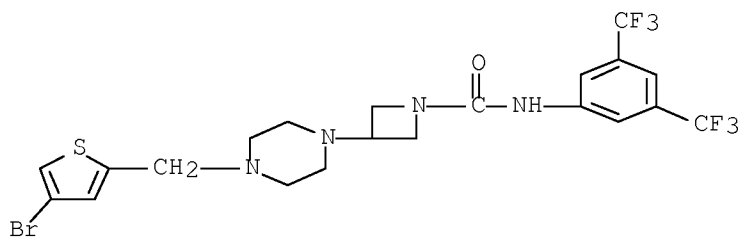
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-methyl-2-thienyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



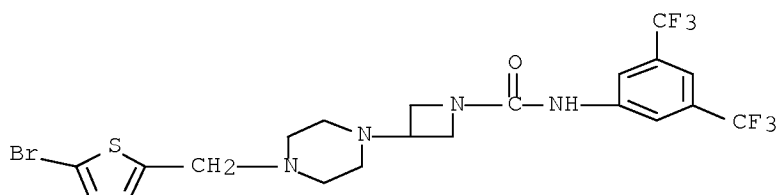
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(4-bromo-2-thienyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



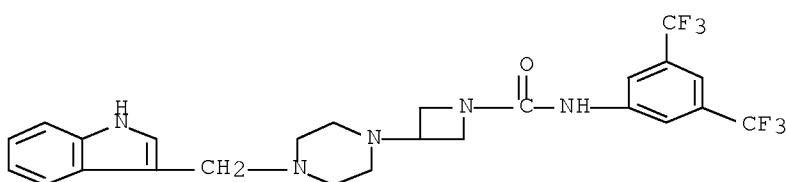
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(5-bromo-2-thienyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



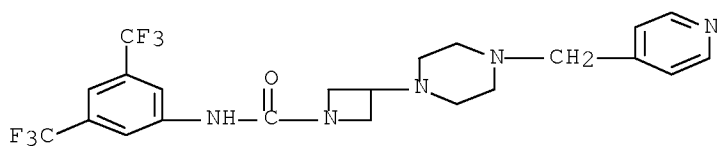
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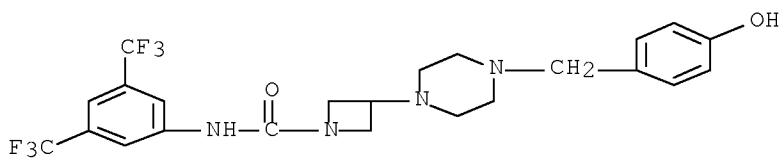
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(4-pyridinylmethyl)-1-piperazinyl]- (CA INDEX NAME)



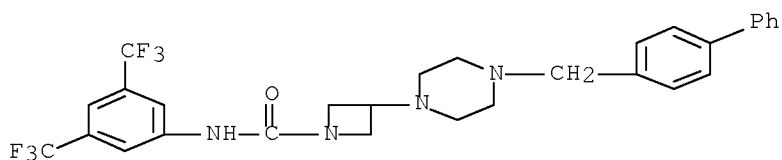
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CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(4-hydroxyphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



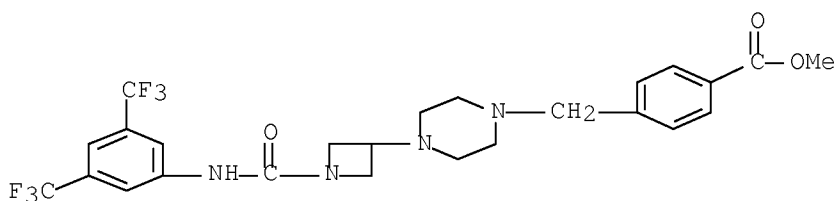
RN 858657-27-9 CAPLUS

CN 1-Azetidinecarboxamide, 3-[4-([1,1'-biphenyl]-4-ylmethyl)-1-piperazinyl]-N-[3,5-bis(trifluoromethyl)phenyl]- (CA INDEX NAME)



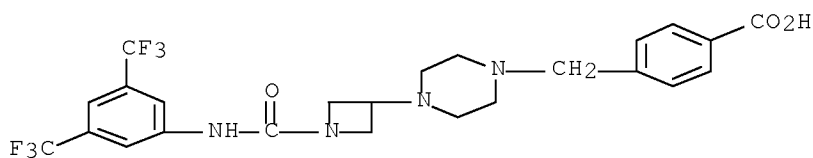
RN 858657-28-0 CAPLUS

CN Benzoic acid, 4-[[4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]-1-piperazinyl]methyl]-, methyl ester (CA INDEX NAME)



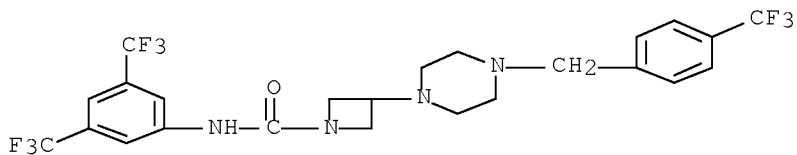
RN 858657-29-1 CAPLUS

CN Benzoic acid, 4-[[4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]-1-piperazinyl]methyl]- (CA INDEX NAME)



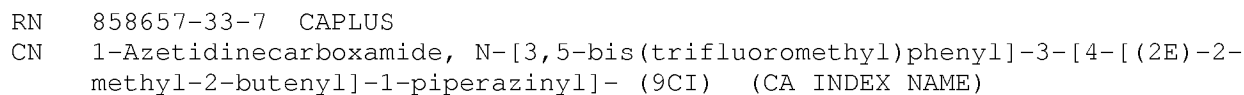
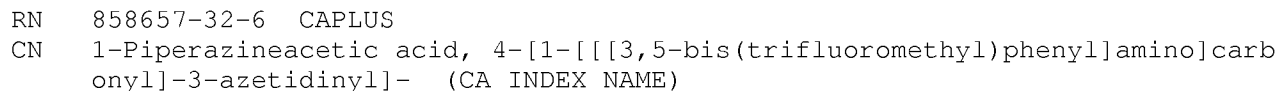
RN 858657-30-4 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(trifluoromethyl)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)

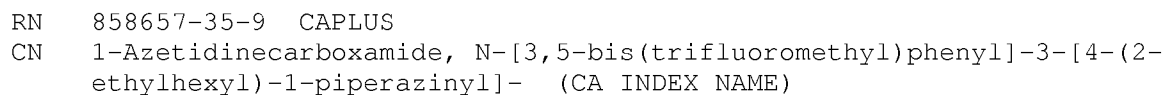


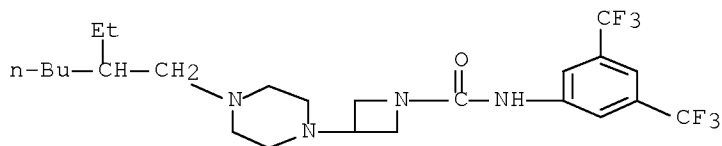
RN 858657-31-5 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(4-methylphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



RN	858657-34-8	CAPLUS
CN	1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(2-methylpropyl)-1-piperazinyl]- (CA INDEX NAME)	

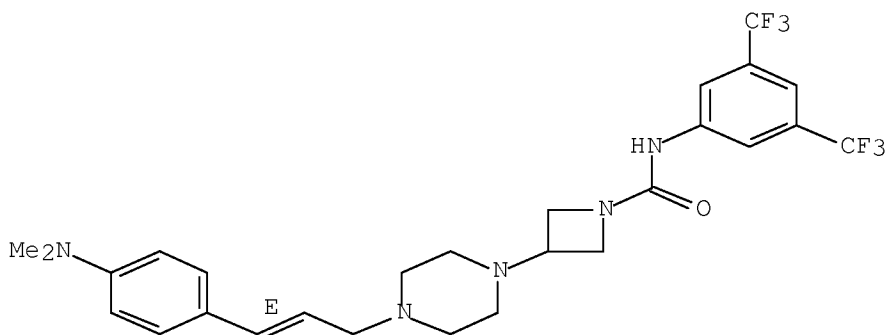




RN 858657-36-0 CAPLUS

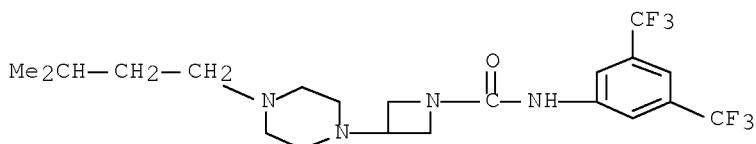
CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(2E)-3-[4-(dimethylamino)phenyl]-2-propenyl]-1-piperazinyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



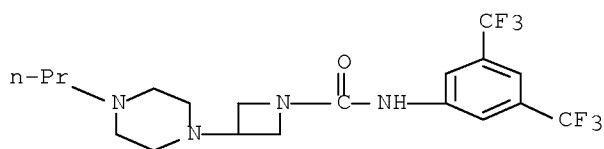
RN 858657-37-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(3-methylbutyl)-1-piperazinyl]- (CA INDEX NAME)



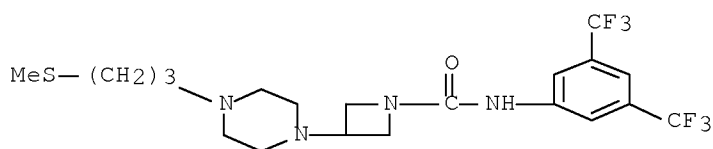
RN 858657-38-2 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(4-propyl-1-piperazinyl)- (CA INDEX NAME)



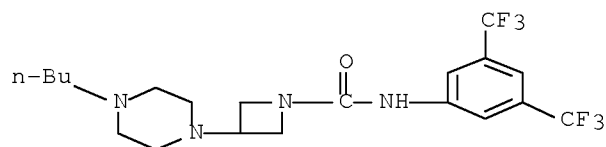
RN 858657-39-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[3-(methylthio)propyl]-1-piperazinyl]- (CA INDEX NAME)



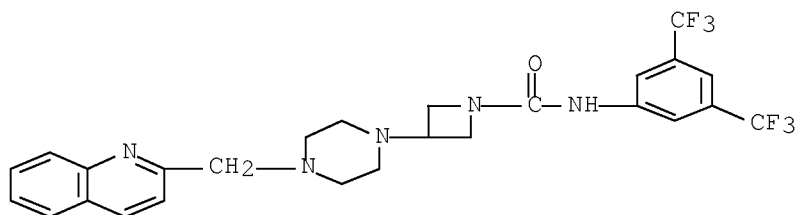
RN 858657-40-6 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-(4-butyl-1-piperazinyl)- (CA INDEX NAME)



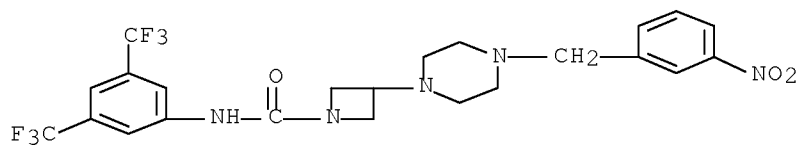
RN 858657-41-7 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(2-quinolinylmethyl)-1-piperazinyl]- (CA INDEX NAME)



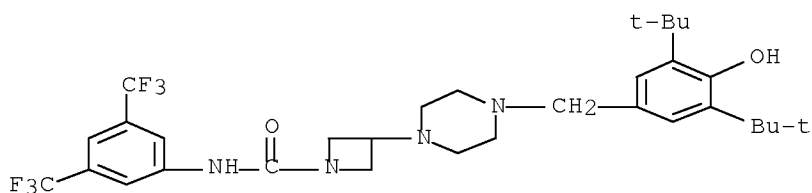
RN 858657-42-8 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-nitrophenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



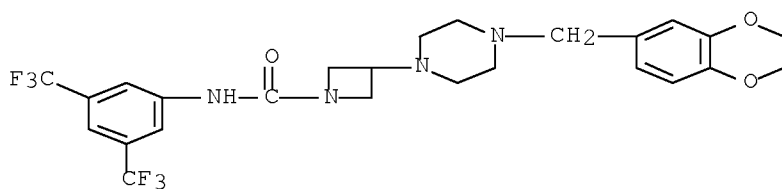
RN 858657-43-9 CAPLUS

CN 1-Azetidinecarboxamide, 3-[4-[[[3,5-bis(1,1-dimethylethyl)-4-hydroxyphenyl]methyl]-1-piperazinyl]-N-[3,5-bis(trifluoromethyl)phenyl]- (CA INDEX NAME)



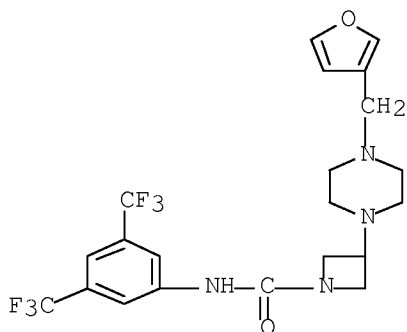
RN 858657-44-0 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(2,3-dihydro-1,4-benzodioxin-6-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



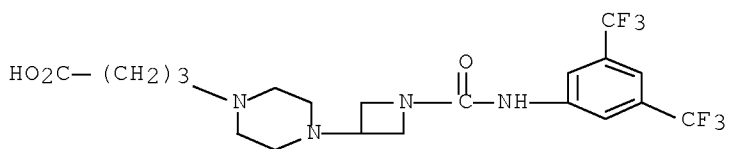
RN 858657-45-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(3-furanylmethyl)-1-piperazinyl]- (CA INDEX NAME)



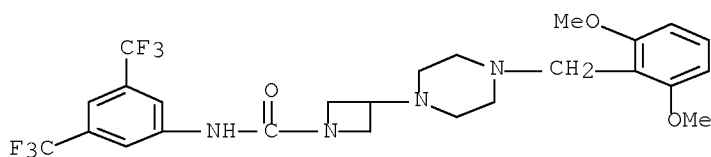
RN 858657-46-2 CAPLUS

CN 1-Piperazinebutanoic acid, 4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]- (CA INDEX NAME)



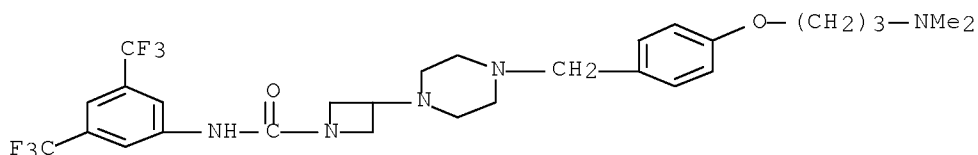
RN 858657-47-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(2,6-dimethoxyphenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



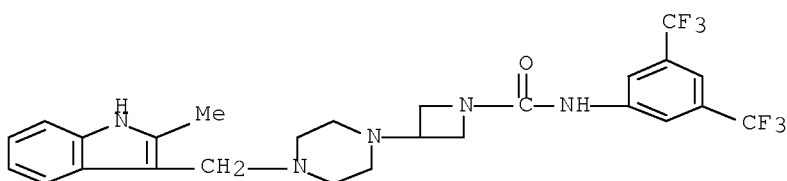
RN 858657-48-4 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-[3-(dimethylamino)propoxy]phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



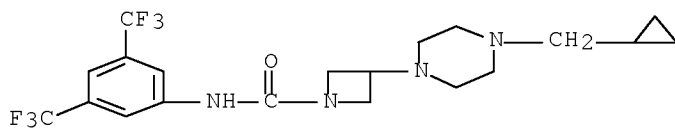
RN 858657-49-5 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(2-methyl-1H-indol-3-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



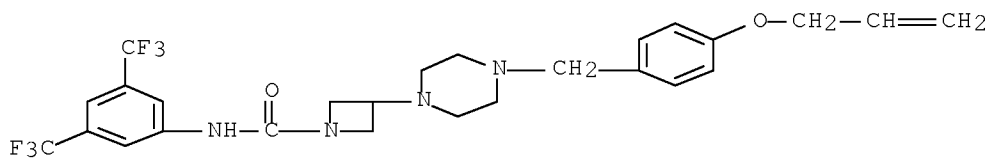
RN 858657-50-8 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(cyclopropylmethyl)-1-piperazinyl]- (CA INDEX NAME)



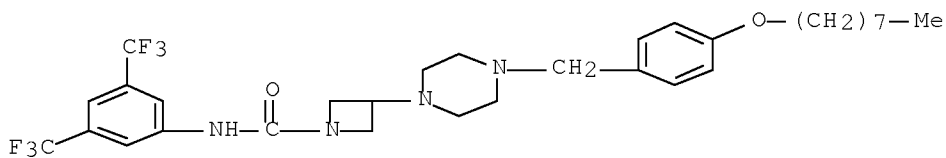
RN 858657-51-9 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(2-propenyloxy)phenyl]methyl]-1-piperazinyl]- (9CI) (CA INDEX NAME)



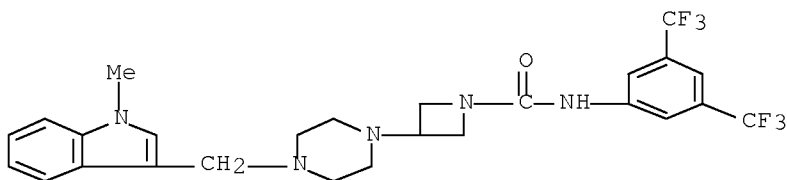
RN 858657-52-0 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(octyloxy)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



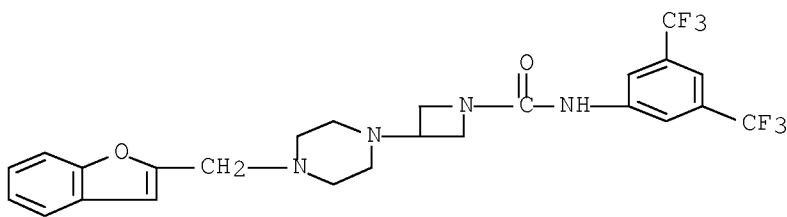
RN 858657-53-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(1-methyl-1H-indol-3-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



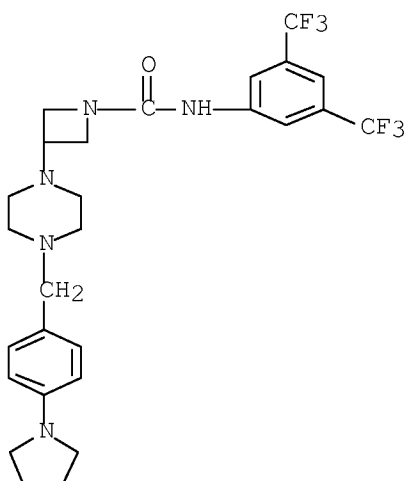
RN 858657-54-2 CAPLUS

CN 1-Azetidinecarboxamide, 3-[4-(2-benzofuranylmethyl)-1-piperazinyl]-N-[3,5-bis(trifluoromethyl)phenyl]- (CA INDEX NAME)



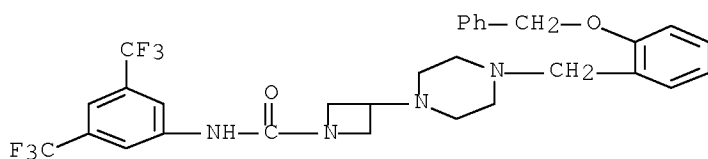
RN 858657-55-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(1-pyrrolidinyl)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



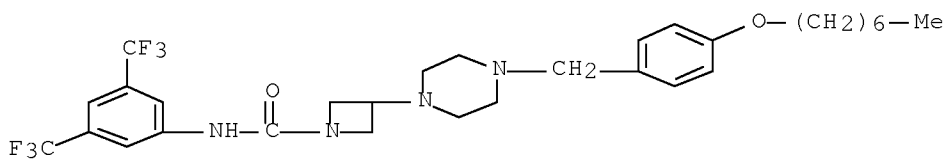
RN 858657-56-4 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[2-(phenylmethoxy)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



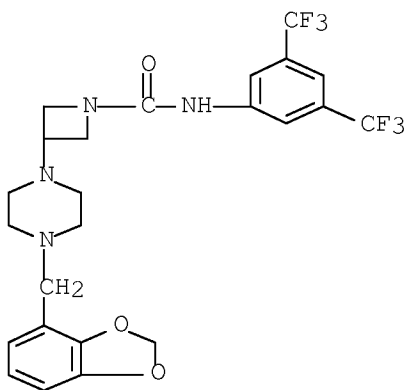
RN 858657-57-5 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(heptyloxy)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



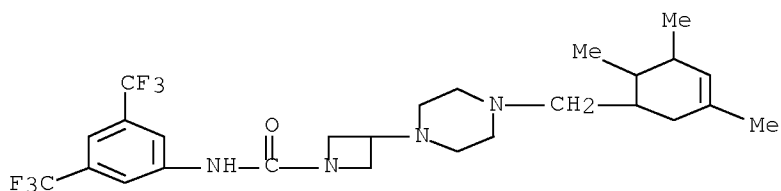
RN 858657-58-6 CAPLUS

CN 1-Azetidinecarboxamide, 3-[4-(1,3-benzodioxol-4-ylmethyl)-1-piperazinyl]-N-[3,5-bis(trifluoromethyl)phenyl]- (CA INDEX NAME)



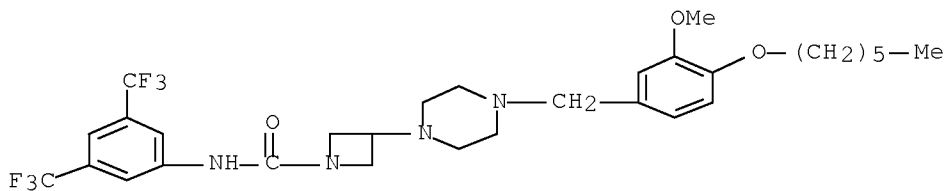
RN 858657-59-7 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3,5,6-trimethyl-3-cyclohexen-1-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



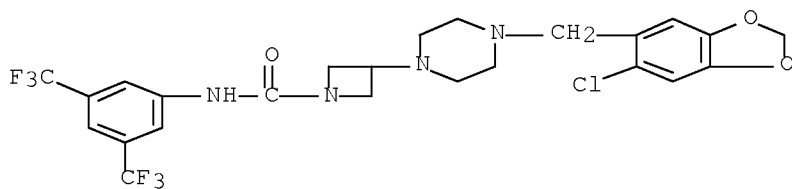
RN 858657-60-0 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(hexyloxy)-3-methoxyphenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



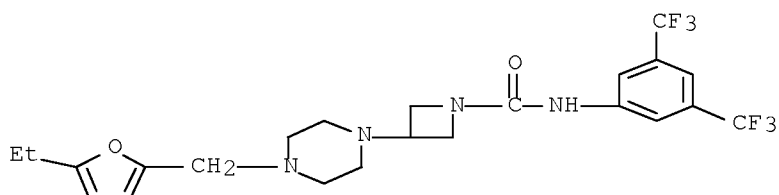
RN 858657-61-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(6-chloro-1,3-benzodioxol-5-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



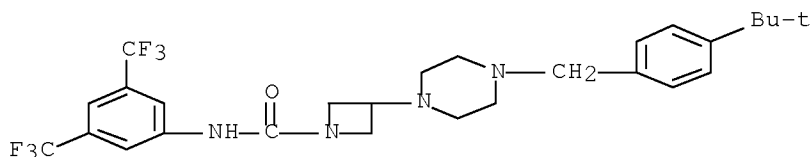
RN 858657-62-2 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(5-ethyl-2-furanyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



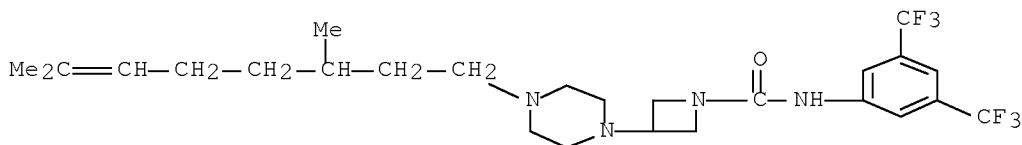
RN 858657-63-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(1,1-dimethylethyl)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



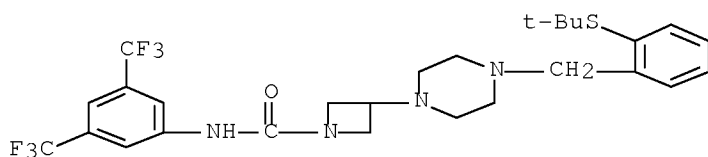
RN 858657-64-4 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(3,7-dimethyl-6-octenyl)-1-piperazinyl]- (9CI) (CA INDEX NAME)



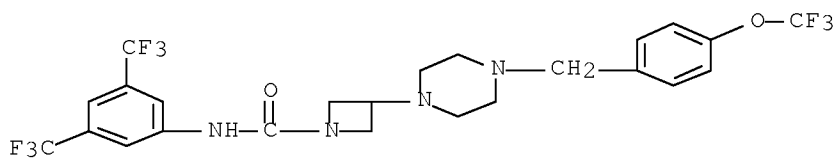
RN 858657-65-5 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[2-[(1,1-dimethylethyl)thio]phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



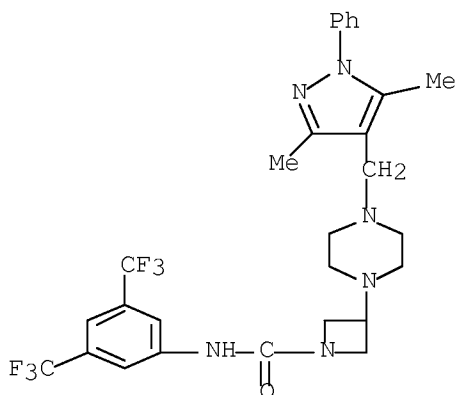
RN 858657-66-6 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(trifluoromethoxy)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



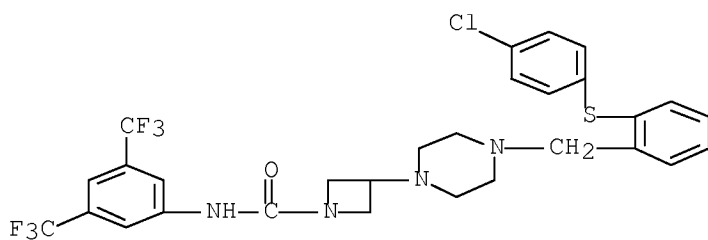
RN 858657-67-7 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3,5-dimethyl-1-phenyl-1H-pyrazol-4-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



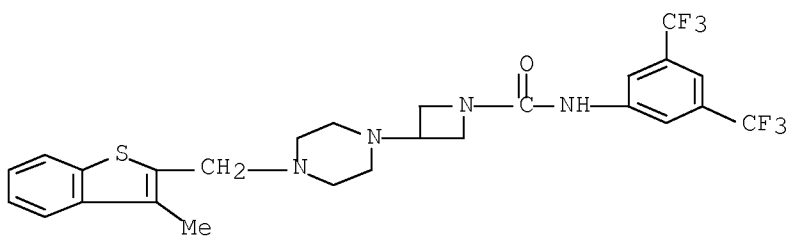
RN 858657-68-8 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[2-[(4-chlorophenyl)thio]phenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



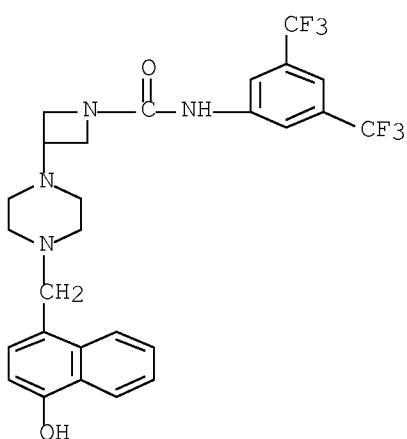
RN 858657-69-9 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-methylbenzo[b]thien-2-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



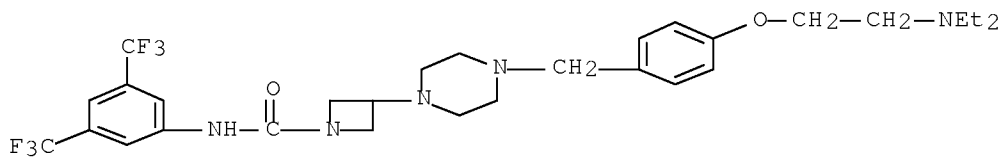
RN 858657-71-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(4-hydroxy-1-naphthalenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



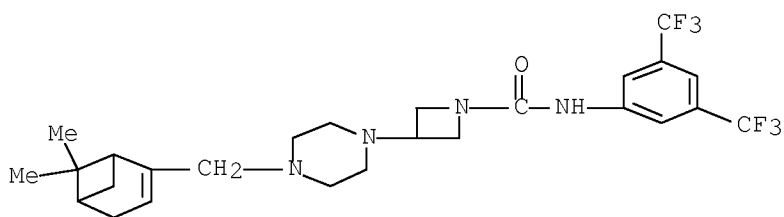
RN 858657-72-4 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-[2-(diethylamino)ethoxy]phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



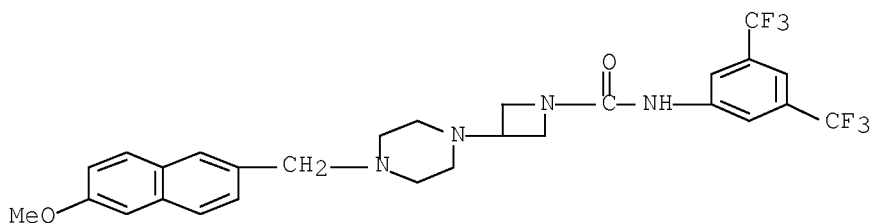
RN 858657-73-5 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(6,6-dimethylbicyclo[3.1.1]hept-2-en-2-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



RN 858657-74-6 CAPLUS

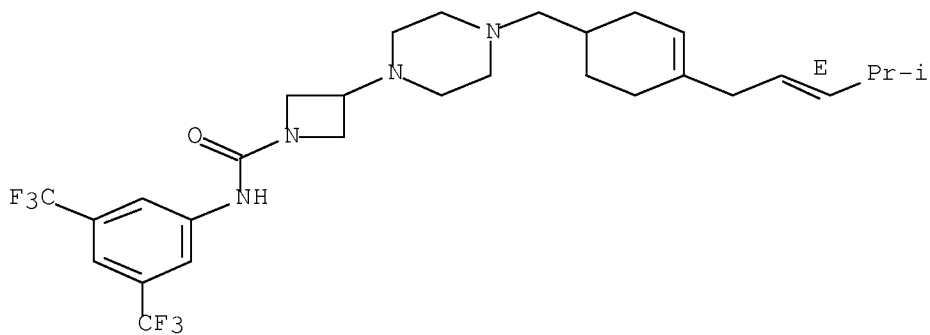
CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(6-methoxy-2-naphthalenyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



RN 858657-75-7 CAPLUS

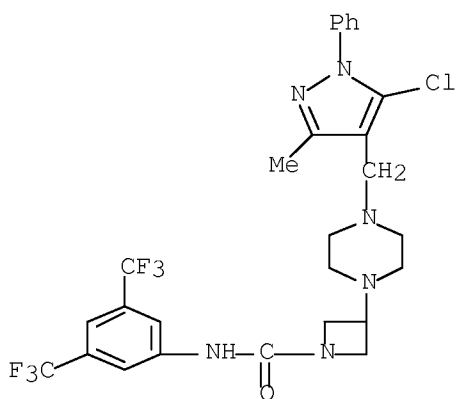
CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-[(2E)-4-methyl-2-pentenyl]-3-cyclohexen-1-yl)methyl]-1-piperazinyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



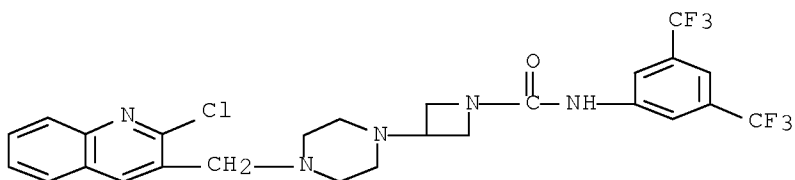
RN 858657-76-8 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(5-chloro-3-methyl-1-phenyl-1H-pyrazol-4-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



RN 858657-77-9 CAPLUS

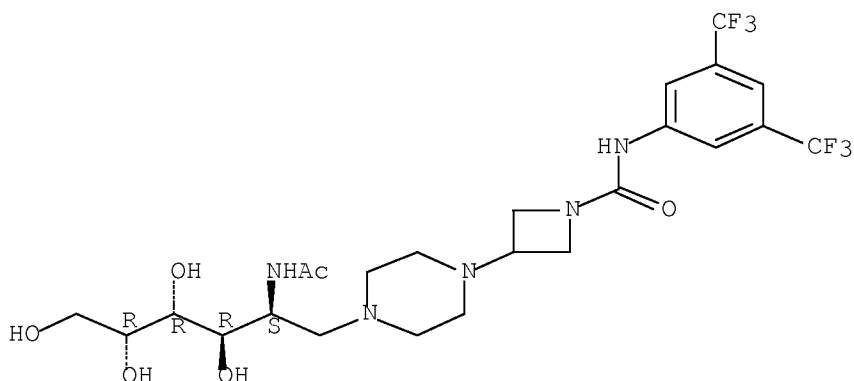
CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(2-chloro-3-quinolinyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



RN 858657-78-0 CAPLUS

CN D-Galactitol, 2-(acetamido)-1-[4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]-1-piperazinyl]-1,2-dideoxy- (CA INDEX NAME)

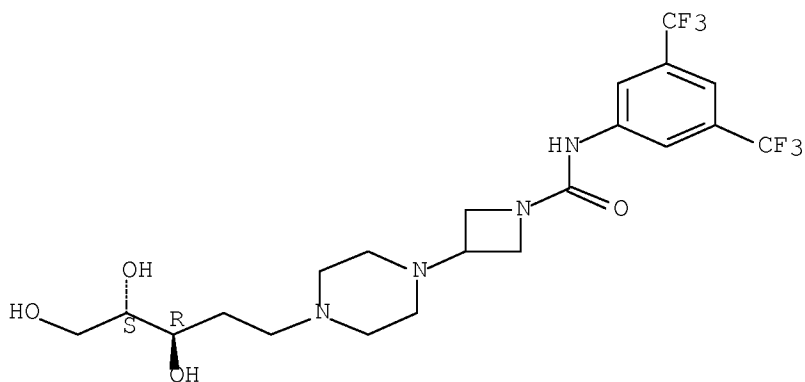
Absolute stereochemistry.



RN 858657-79-1 CAPLUS

CN D-erythro-Pentitol, 5-[4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]-1-piperazinyl]-4,5-dideoxy- (CA INDEX NAME)

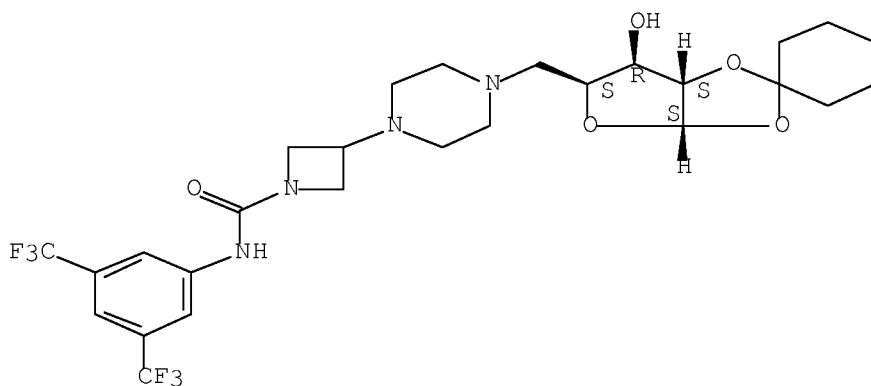
Absolute stereochemistry.



RN 858657-80-4 CAPLUS

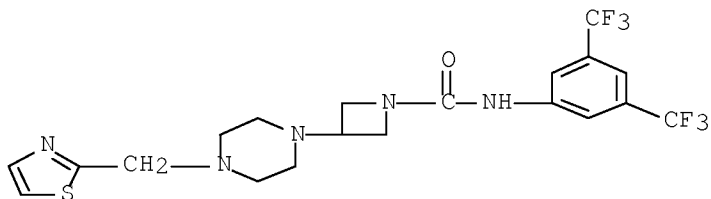
CN α -L-Xylofuranose, 5-[4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]-1-piperazinyl]-1,2-O-cyclohexylidene-5-deoxy- (CA INDEX NAME)

Absolute stereochemistry.



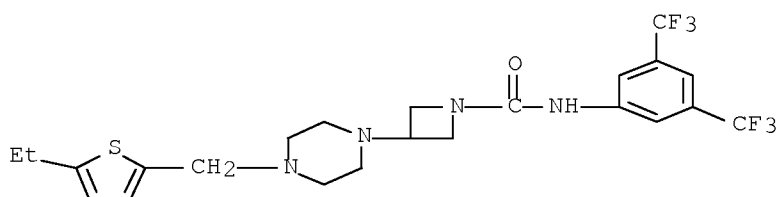
RN 858657-81-5 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-(2-thiazolylmethyl)-1-piperazinyl]- (CA INDEX NAME)



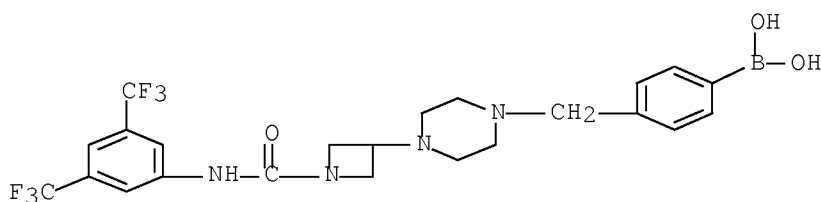
RN 858657-82-6 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(5-ethyl-2-thienyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



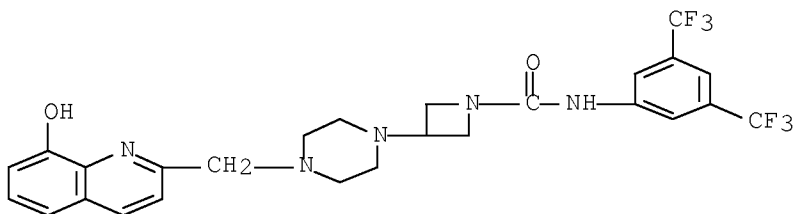
RN 858657-83-7 CAPLUS

CN Boronic acid, [4-[[4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]]-1-piperazinyl]methyl]phenyl]- (9CI) (CA INDEX NAME)



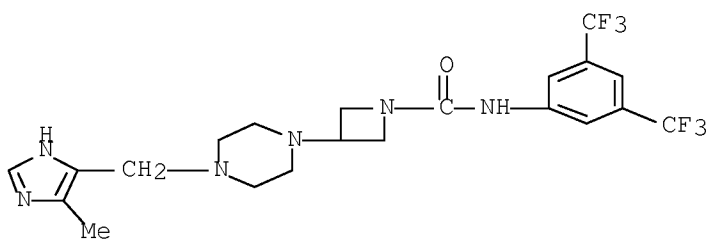
RN 858657-84-8 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(8-hydroxy-2-quinolinyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



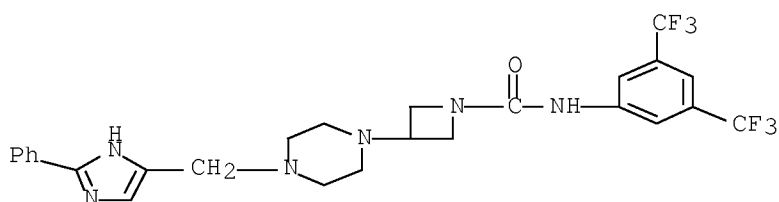
RN 858657-85-9 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(5-methyl-1H-imidazol-4-yl)methyl]-1-piperazinyl]- (9CI) (CA INDEX NAME)



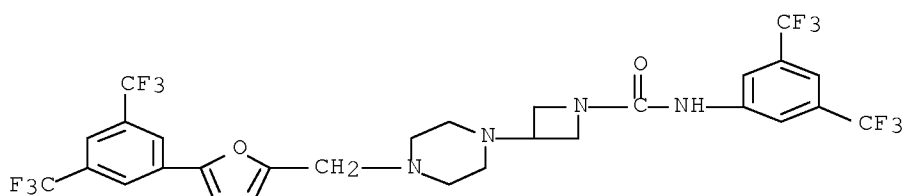
RN 858657-86-0 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(2-phenyl-1H-imidazol-4-yl)methyl]-1-piperazinyl]- (9CI) (CA INDEX NAME)



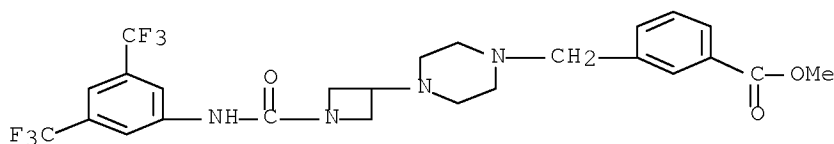
RN 858657-87-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[5-[3,5-bis(trifluoromethyl)phenyl]-2-furanyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



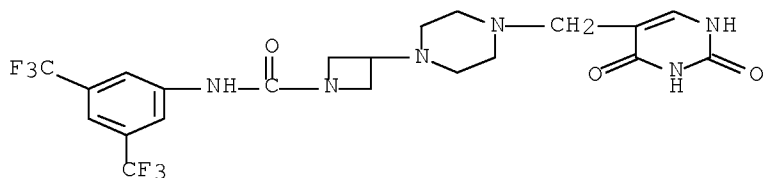
RN 858657-88-2 CAPLUS

CN Benzoic acid, 3-[[4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]-1-piperazinyl]methyl]-, methyl ester (CA INDEX NAME)



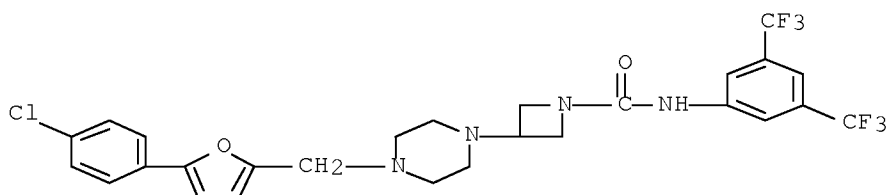
RN 858657-89-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(1,2,3,4-tetrahydro-2,4-dioxo-5-pyrimidinyl)methyl]-1-piperazinyl]- (CA INDEX NAME)



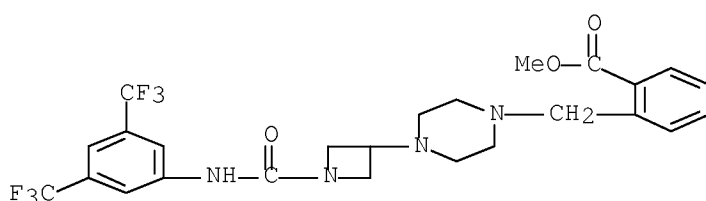
RN 858657-90-6 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[5-(4-chlorophenyl)-2-furanyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



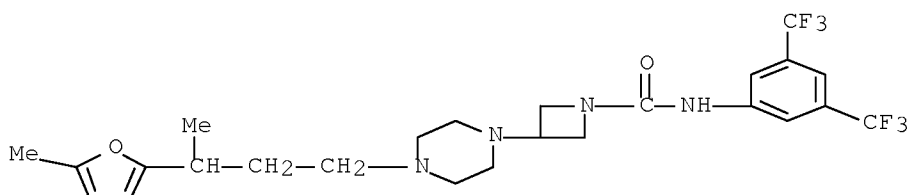
RN 858657-91-7 CAPLUS

CN Benzoic acid, 2-[[4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]-1-piperazinyl]methyl]-, methyl ester (CA INDEX NAME)



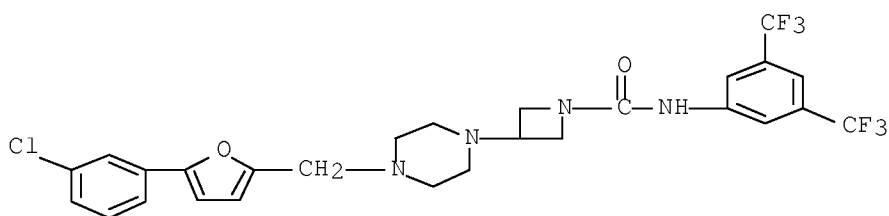
RN 858657-92-8 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[3-(5-methyl-2-furanyl)butyl]-1-piperazinyl]- (CA INDEX NAME)



RN 858657-93-9 CAPLUS

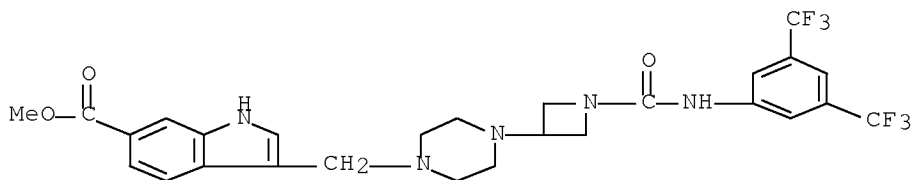
CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[5-(3-chlorophenyl)-2-furanyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



RN 858657-94-0 CAPLUS

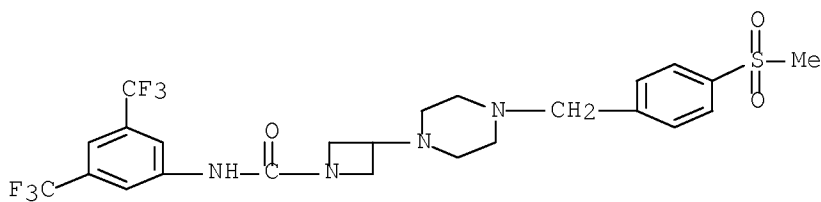
CN 1H-Indole-6-carboxylic acid, 3-[[4-[1-[[[3,5-bis(trifluoromethyl)phenyl]amino]carbonyl]-3-azetidiny]-1-piperazinyl]methyl]-, methyl ester (CA INDEX NAME)

INDEX NAME)



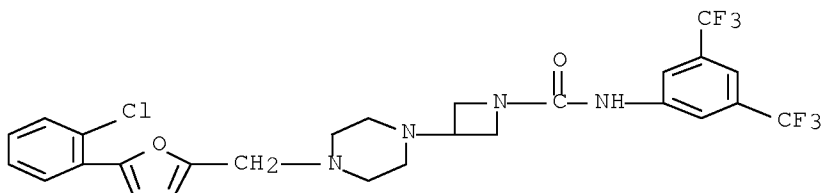
RN 858657-95-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[4-(methylsulfonyl)phenyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



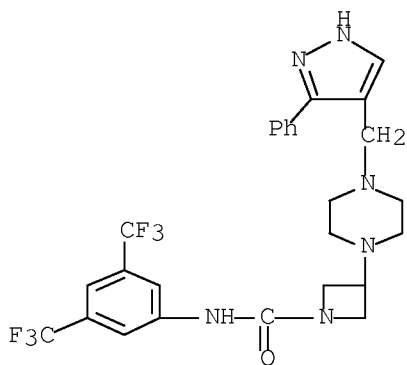
RN 858657-96-2 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[5-(2-chlorophenyl)-2-furanyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



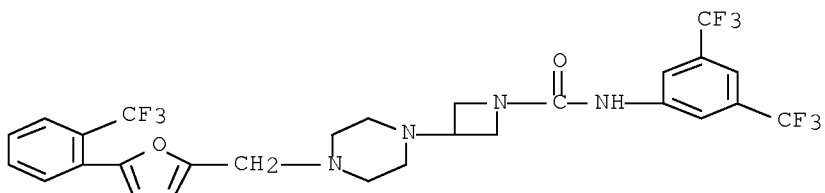
RN 858657-97-3 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[(3-phenyl-1H-pyrazol-4-yl)methyl]-1-piperazinyl]- (CA INDEX NAME)



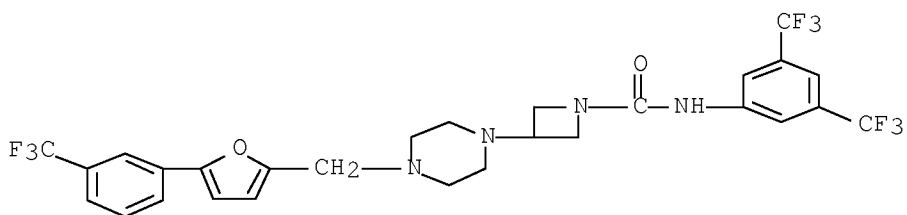
RN 858657-98-4 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[5-[2-(trifluoromethyl)phenyl]-2-furanyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



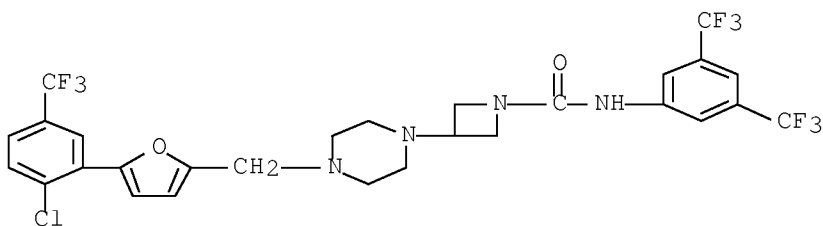
RN 858657-99-5 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[5-[3-(trifluoromethyl)phenyl]-2-furanyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



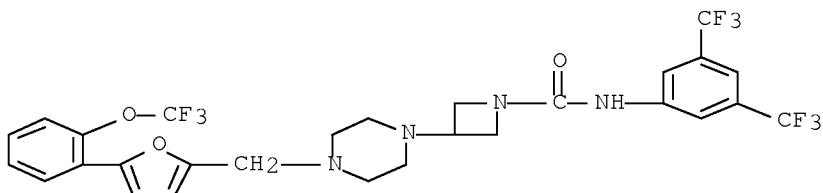
RN 858658-00-1 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[5-[2-chloro-5-(trifluoromethyl)phenyl]-2-furanyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



RN 858658-01-2 CAPLUS

CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[[5-[2-(trifluoromethoxy)phenyl]-2-furanyl]methyl]-1-piperazinyl]- (CA INDEX NAME)



REFERENCE COUNT: 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 2 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2005:490293 CAPLUS Full-text

DOCUMENT NUMBER: 143:43903

TITLE: Preparation of piperazinylguanidinoquinazolinones as melanocortin-4 receptor (MCR-4) agonists with reduced bioaccumulation

INVENTOR(S): Boyce, Rustum S.; Speake, Jason D.; Phillips, James

PATENT ASSIGNEE(S): Chiron Corporation, USA; Glaxosmithkline

SOURCE: PCT Int. Appl., 199 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

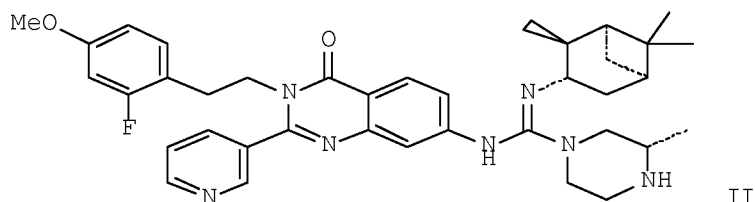
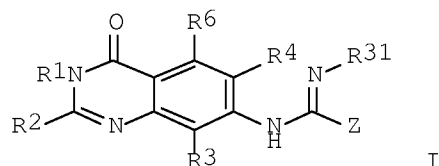
LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005051391	A1	20050609	WO 2004-US39020	20041119 <--
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
RW:	BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
AU 2004293012	A1	20050609	AU 2004-293012	20041119 <--
CA 2545601	A1	20050609	CA 2004-2545601	20041119 <--

US 20050192297	A1	20050901	US 2004-993147	20041119 <--
EP 1686996	A1	20060809	EP 2004-811698	20041119 <--
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK, IS				
CN 1901916	A	20070124	CN 2004-80039762	20041119 <--
JP 2007511612	T	20070510	JP 2006-541565	20041119 <--
MX 2006PA05736	A	20061214	MX 2006-PA5736	20060519 <--
IN 2006KN01610	A	20070511	IN 2006-KN1610	20060612 <--
PRIORITY APPLN. INFO.:			US 2003-523336P	P 20031119 <--
			US 2003-524492P	P 20031124 <--
			WO 2004-US39020	W 20041119
OTHER SOURCE(S): MARPAT 143:43903				
GI				



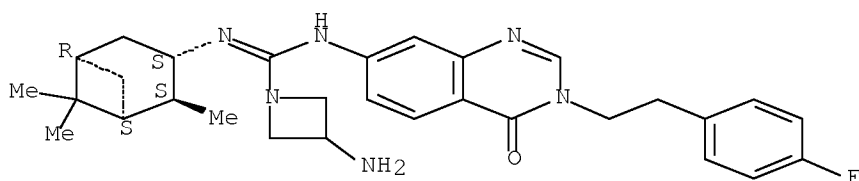
AB Title compds. [I; R1 = (substituted) aralkyl, heteroarylalkyl, aryl, heteroaryl, cycloalkyl, heterocycloalkyl, cycloalkylalkyl, alkenyl, alkynyl, alkyl; R2 = H, (substituted) aralkyl, heteroarylalkyl, alkoxy, alkylamino, dialkylamino, aryl, heteroaryl, heterocyclyl, cycloalkyl, heterocycloalkyl, cycloalkylalkyl, alkenyl, alkynyl, alkyl; R3, R4, R6 = H, Cl, F, Br, iodo, OH, NH2, cyano, NO2, (substituted) alkoxy, alkyl; R31 = H, (substituted) alkyl, aryl, alkenyl, alkynyl, cycloalkyl, heteroaryl, heterocyclyl, heterocyclylalkyl, aralkyl, heteroarylalkyl, cycloalkylalkyl; Z = (substituted) 3-oxopiperazinyl; and tautomers], were prepared Thus, title compound (II) (preparation via coupling of 6-methylpiperazin-2-one with the corresponding quinazolinylthiourea derivative in the presence of polymer-supported carbodiimide) showed a plasma half life of 1.9 h in mice.

IT 817622-71-2P 817627-62-6P 817627-64-8P
 RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
 (preparation of piperazinylguanidinoquinazolinones as melanocortin-4 receptor (MCR-4) agonists with reduced bioaccumulation)

RN 817622-71-2 CAPLUS

CN 1-Azetidinecarboximidamide, 3-amino-N-[3-[2-(4-fluorophenyl)ethyl]-3,4-dihydro-4-oxo-7-quinazolinyl]-N'-[(1S,2S,3S,5R)-2,6,6-trimethylbicyclo[3.1.1]hept-3-yl]- (CA INDEX NAME)

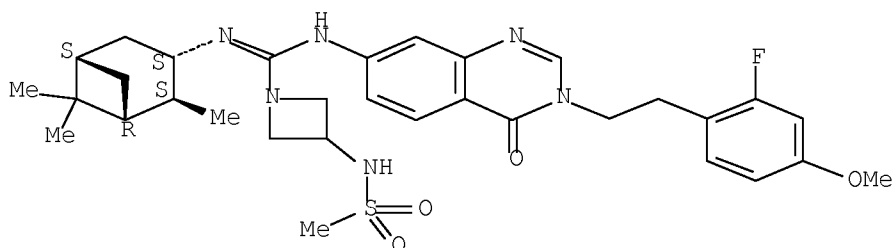
Absolute stereochemistry.



RN 817627-62-6 CAPLUS

CN 1-Azetidinecarboximidamide, N-[3-[2-(2-fluoro-4-methoxyphenyl)ethyl]-3,4-dihydro-4-oxo-7-quinazolinyl]-3-[(methylsulfonyl)amino]-N'-[(1R,2S,3S,5S)-2,6,6-trimethylbicyclo[3.1.1]hept-3-yl]- (CA INDEX NAME)

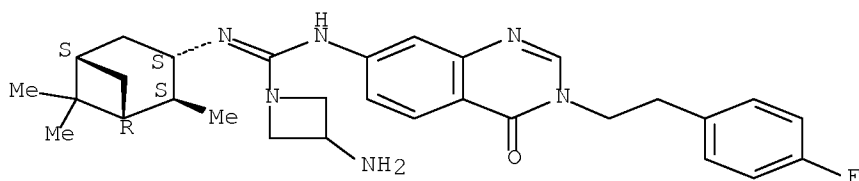
Absolute stereochemistry.



RN 817627-64-8 CAPLUS

CN 1-Azetidinecarboximidamide, 3-amino-N-[3-[2-(4-fluorophenyl)ethyl]-3,4-dihydro-4-oxo-7-quinazolinyl]-N'-[(1R,2S,3S,5S)-2,6,6-trimethylbicyclo[3.1.1]hept-3-yl]- (CA INDEX NAME)

Absolute stereochemistry.



IT 819051-14-4

RL: RCT (Reactant); RACT (Reactant or reagent)

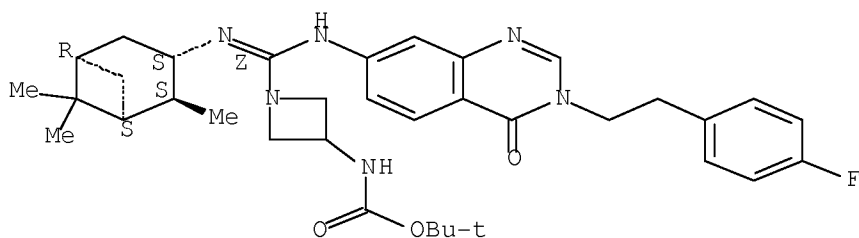
(preparation of piperazinyguanidinoquinazolinones as melanocortin-4 receptor (MCR-4) agonists with reduced bioaccumulation)

RN 819051-14-4 CAPLUS

CN Carbamic acid, [1-[(Z)-[[3-[2-(4-fluorophenyl)ethyl]-3,4-dihydro-4-oxo-7-quinazolinyl]amino][[(1S,2S,3S,5R)-2,6,6-trimethylbicyclo[3.1.1]hept-3-yl]imino]methyl]-3-azetidiny]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Double bond geometry as shown.



REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 3 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN
ACCESSION NUMBER: 2004:1156498 CAPLUS Full-text

DOCUMENT NUMBER: 142:93848

TITLE: Preparation of guanidino-substituted quinazolinone compounds as MC4-R agonists

INVENTOR(S): Boyce, Rustum S.; Aurrecoechea, Natalia; Chu, Daniel; Smith, Aaron; Conlee, Christopher R.; Thompson, Brian D.; De Armas, Kuntz Judith; Musso, David L.; Barvian, Kevin K.; Thomson, Stephen A.; Swain, William R.; Du, Kien S.; Chauder, Brian A.; Speake, Jason D.; Bishop, Michael J.

PATENT ASSIGNEE(S): Chiron Corporation, USA; Glaxosmithkline

SOURCE: PCT Int. Appl., 277 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

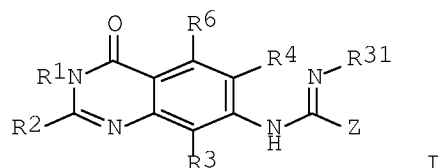
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004112793	A1	20041229	WO 2004-US15959	20040521 <--
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
AU 2004249120	A1	20041229	AU 2004-249120	20040521 <--
CA 2523015	A1	20041229	CA 2004-2523015	20040521 <--
US 20050059662	A1	20050317	US 2004-850967	20040521 <--
EP 1651229	A1	20060503	EP 2004-776069	20040521 <--
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK				
CN 1829517	A	20060906	CN 2004-80013951	20040521 <--
JP 2007501861	T	20070201	JP 2006-533275	20040521 <--
IN 2005KN02103	A	20070720	IN 2005-KN2103	20051024 <--
MX 2005PA12483	A	20060929	MX 2005-PA12483	20051118 <--
PRIORITY APPLN. INFO.:			US 2003-473317P	P 20030523 <--
			US 2003-523336P	P 20031119 <--

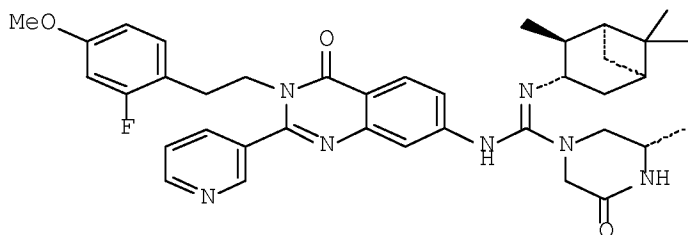
US 2003-524492P
WO 2004-US15959

P 20031124 <--
W 20040521

OTHER SOURCE(S): MARPAT 142:93848
GI



I



II

AB A variety of small mol., guanidine-containing mols. capable of acting as MC4-R agonists such as I-III [Z1 = CR4, N; Z2 = CR5, N; Z3 = CR6, N; R1 = (un)substituted arylalkyl, heteroarylalkyl, aryl, heteroaryl, etc.; R2 = H, alkyl, aryl, etc.; R3 = H, arylalkyl, aryl, etc.; R4-R6 = H, Cl, I, F, Br, OH, etc.; W = IV (wherein R11, R12 = H, (un)substituted alkyl, aryl, etc.; at least one of R11 and R12 is (un)substituted heterocyclalkyl; R13 = H, (un)substituted aryl, alkyl, etc.; R14 = H, (un)substituted alkyl, cycloalkyl, etc.)] are provided. General procedures used in the synthesis of compds. I-III are described. E.g., a multi-step synthesis of (1S,2S,3S,5R)-V, was given. The exemplified compds. I-III were tested against MC4-R and exhibited -logEC50 values above about 3. The compds. I are useful in treating MC4-R mediated diseases such as obesity and type II diabetes. The pharmaceutical composition comprising the compound I is disclosed.

IT 817622-71-2F 817627-62-6F 817627-64-8F

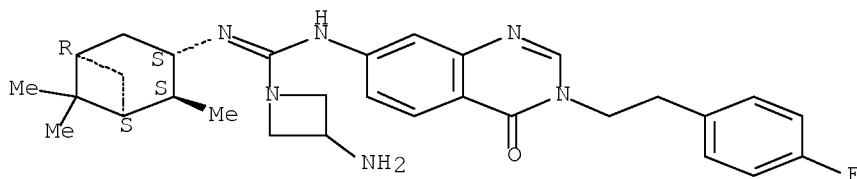
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of guanidino-substituted quinazolinone compds. as MC4-R agonists)

RN 817622-71-2 CAPLUS

CN 1-Azetidinecarboximidamide, 3-amino-N-[3-[2-(4-fluorophenyl)ethyl]-3,4-dihydro-4-oxo-7-quinazolinyl]-N'-[(1S,2S,3S,5R)-2,6,6-trimethylbicyclo[3.1.1]hept-3-yl]- (CA INDEX NAME)

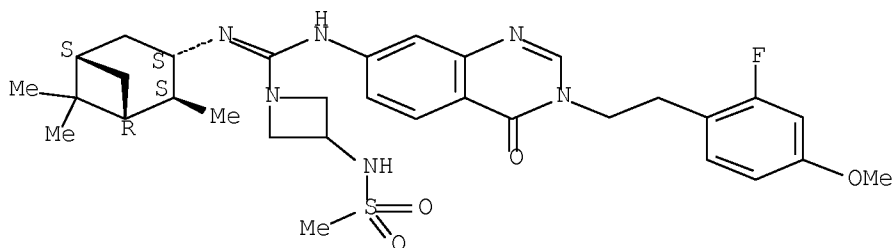
Absolute stereochemistry.



RN 817627-62-6 CAPLUS

CN 1-Azetidinecarboximidamide, N-[3-[2-(2-fluoro-4-methoxyphenyl)ethyl]-3,4-dihydro-4-oxo-7-quinazolinyl]-3-[(methylsulfonyl)amino]-N'-[(1R,2S,3S,5S)-2,6,6-trimethylbicyclo[3.1.1]hept-3-yl]- (CA INDEX NAME)

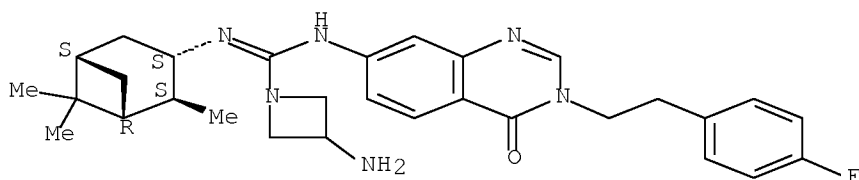
Absolute stereochemistry.



RN 817627-64-8 CAPLUS

CN 1-Azetidinecarboximidamide, 3-amino-N-[3-[2-(4-fluorophenyl)ethyl]-3,4-dihydro-4-oxo-7-quinazolinyl]-N'-[(1R,2S,3S,5S)-2,6,6-trimethylbicyclo[3.1.1]hept-3-yl]- (CA INDEX NAME)

Absolute stereochemistry.



IT 819051-14-4

RL: RCT (Reactant); RACT (Reactant or reagent)

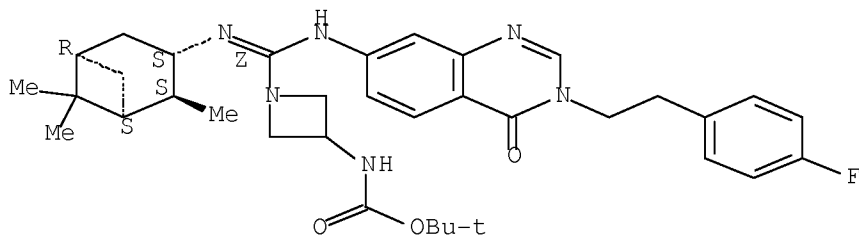
(preparation of guanidino-substituted quinazolinone compds. as MC4-R agonists)

RN 819051-14-4 CAPLUS

CN Carbamic acid, [1-[(Z)-[[3-[2-(4-fluorophenyl)ethyl]-3,4-dihydro-4-oxo-7-quinazolinyl]amino][[(1S,2S,3S,5R)-2,6,6-trimethylbicyclo[3.1.1]hept-3-yl]imino]methyl]-3-azetidiny]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

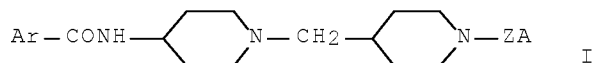
Double bond geometry as shown.



REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 4 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN
ACCESSION NUMBER: 2004:819910 CAPLUS Full-text
DOCUMENT NUMBER: 141:307557
TITLE: 1,4-Disubstituted piperidine derivatives as 5-HT4 agonists, pharmaceutical compns. containing them, and their use
INVENTOR(S): Kato, Shiro; Yamazaki, Hiroshi; Hirokawa, Yoshimi; Suga, Yoko; Yoshida, Naoyuki; Morikage, Yukiko; Morikage, Kazuo; Oku, Seiko
PATENT ASSIGNEE(S): Dainippon Pharmaceutical Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 31 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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JP 2004277320	A	20041007	JP 2003-69337	20030314 <--
PRIORITY APPLN. INFO.:			JP 2003-69337	20030314 <--
OTHER SOURCE(S):	MARPAT	141:307557		
GI				



AB The derivs. are represented by I [Ar = Q, Q1; R1 = halo; R2 = H, lower alkyl; R3 = H, lower alkyl, lower alkanoyl; R4 = H, lower alkyl; R5, R6 = H, lower alkyl; h = 1, 2, 3; Z= CO, CS, SO2; A = NR7R8, Q2; R7 = (un)substituted lower alkyl, lower alkenyl, lower alkynyl, lower alkoxy; R8 = H, lower alkyl, lower alkenyl, lower alkynyl; if R7 = lower alkyl, then R8 is lower alkenyl, lower alkynyl; p + q = 1-7; Q = CH2, CH:CH; R9 = H, (un)substituted lower alkyl, lower alkoxy, OH, amino, mono-or di(lower alkyl)amino; R10 = H, lower alkyl; except cases where q + p = 3-5, Q = CH2, and R9 and R10 = H] or their physiol. acceptable acid salts. Also claimed are pharmaceutical compns., 5-HT4 receptor agonists, and digestive tract motility promoters or digestive tract function improvers containing I or their salts. Thus, 4-amino-5-chloro-2-methoxy-N-[1-[1-(N-methyl-N-2-propynylaminocarbonyl)-4-piperidinylmethyl]-4-piperidinyl]benzamide fumarate (II), prepared from N-methylpropargylamine and 4-amino-5-chloro-2-methoxy-N-[1-(4-piperidinylmethyl)-4-piperidinyl]benzamide, inhibited binding of GR113808 to a guinea pig 5-HT4 receptor preparation at IC50 4.93 nM, vs. 23.0 nM of cisapride. Pharmaceutical preps. containing II were also given.

IT 767306-16-1P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of N-[(piperidinylmethyl)piperidinyl]benzamide derivs. as 5-

TH4

receptor agonists)

RN 767306-16-1 CAPLUS

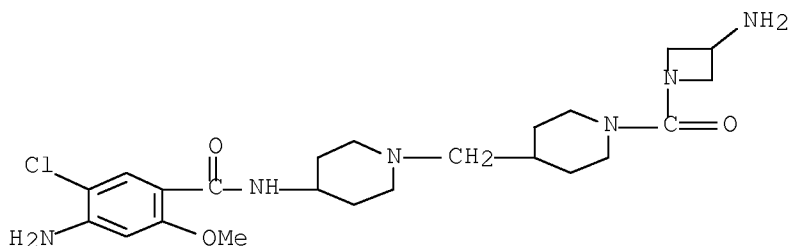
CN Benzamide, 4-amino-N-[1-[1-[(3-amino-1-azetidiny)carbonyl]-4-

piperidinyl)methyl]-4-piperidinyl]-5-chloro-2-methoxy-,
(2E)-2-butenedioate (1:2) (CA INDEX NAME)

CM 1

CRN 767306-15-0

CMF C23 H35 Cl N6 O3

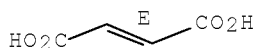


CM 2

CRN 110-17-8

CMF C4 H4 O4

Double bond geometry as shown.



L17 ANSWER 5 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN
ACCESSION NUMBER: 2004:390255 CAPLUS [Full-text](#)
DOCUMENT NUMBER: 140:406684
TITLE: Synthesis of (thiadiazolyliminoacetamido)(pyrazoliummet
hyl)cephem compounds as antimicrobial agents
INVENTOR(S): Ohki, Hidenori; Okuda, Shinya; Yamanaka, Toshio;
Ohgaki, Masaru; Toda, Ayako; Kawabata, Kohji; Inoue,
Satoshi; Misumi, Keiji; Itoh, Kenji; Satoh, Kenji
PATENT ASSIGNEE(S): Fujisawa Pharmaceutical Co., Ltd., Japan; Wakunaga
Pharmaceutical Co., Ltd.; et al.
SOURCE: PCT Int. Appl., 137 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004039814	A1	20040513	WO 2003-JP13684	20031027 <--
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN,				

TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY,
KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES,
FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR,
BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

CA 2504730	A1	20040513	CA 2003-2504730	20031027 <--
AU 2003274754	A1	20040525	AU 2003-274754	20031027 <--
EP 1556389	A1	20050727	EP 2003-758919	20031027 <--
EP 1556389	B1	20070725		

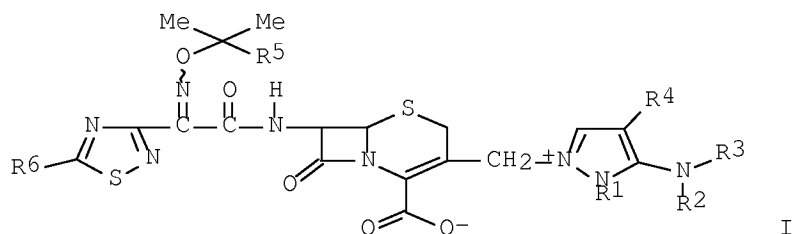
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK

BR 2003015188	A	20050906	BR 2003-15188	20031027 <--
CN 1708502	A	20051214	CN 2003-80102642	20031027 <--
JP 2006506459	T	20060223	JP 2005-501847	20031027 <--
AT 368042	T	20070815	AT 2003-758919	20031027 <--
ES 2290498	T3	20080216	ES 2003-758919	20031027 <--
US 20040132994	A1	20040708	US 2003-695895	20031030 <--
US 7129232	B2	20061031		
IN 2005KN00936	A	20060811	IN 2005-KN936	20050519 <--
ZA 2005004181	A	20060830	ZA 2005-4181	20050523 <--
US 20070037786	A1	20070215	US 2006-585262	20061024 <--

PRIORITY APPLN. INFO.:

			AU 2002-952355	A 20021030 <--
			AU 2003-904813	A 20030904 <--
			WO 2003-JP13684	W 20031027 <--
			US 2003-695895	A1 20031030 <--

OTHER SOURCE(S): MARPAT 140:406684
GI



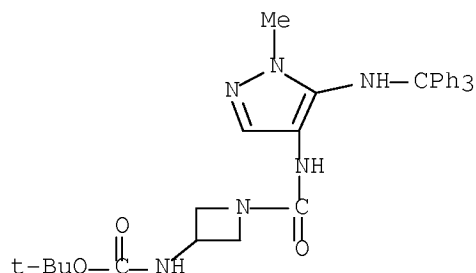
AB Cephem derivs. I [R1 = (hydroxy/halo)alkyl; R2 = H, amino protecting group; R1R2 = alkylene, alkenylene; R3 = H, alkyl; R4 = N(R7) (A)k (NH)mOn (CHR8)q (CH2)pR9, A = C:X, COCO, COCH2CO, etc., R7 = H, alkyl, amino protecting group, R8 = H, OH, R9 = amino, dialkylamino, protected amino, etc., k, m, n, q = independently 0, 1, p = 0-3, X = O, NH; R5 = carboxy, protected carboxy; R6 = amino, protected amino] were prepared to be used as antimicrobial agents. Thus, benzyhydriyl 7β-[(Z)-2-(5-tert-butoxycarbonylamino-1,2,4-thiadiazol-3-yl)-2-(1-tert-butoxycarbonyl-1-methylethoxyimino)acetamido]-3-chloromethyl-3-cephem-4-carboxylate reacted with 5-amino-4-(3-(2-[(tert-butoxycarbonyl)amino]ethyl)ureido)-1-methylpyrazole to give 7β-[(Z)-2-(5-amino-1,2,4-thiadiazol-3-yl)-2-(1-carboxy-1-methylethoxyimino)acetamido]-3-[3-amino-4-[3-(2-aminoethyl)ureido]-2-methyl-1-pyrazolio]methyl-3-cephem-4-carboxylate. The prepared cepheims were tested in vitro for antibacterial activity against Pseudomonas aeruginosa FP 1380.

IT 689294-51-7F
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(intermediate; synthesis of (thiadiazolyliminoacetamido)(pyrazolomethyl)cephem compds. as antimicrobial agents)

RN 689294-51-7 CAPLUS

CN Carbamic acid, [1-[[[1-methyl-5-[(triphenylmethyl)amino]-1H-pyrazol-4-yl]amino]carbonyl]-3-azetidiny]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



IT 689294-52-8P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

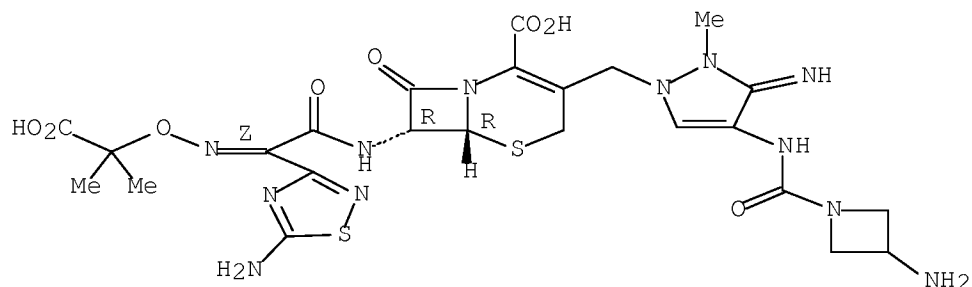
(product; synthesis of (thiadiazolyliminoacetamido)(pyrazolomethyl)cephem compds. as antimicrobial agents)

RN 689294-52-8 CAPLUS

CN 5-Thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylic acid, 3-[[4-[[[(3-amino-1-azetidiny)carbonyl]amino]-2,3-dihydro-3-imino-2-methyl-1H-pyrazol-1-yl]methyl]-7-[[[(2Z)-(5-amino-1,2,4-thiadiazol-3-yl)][(1-carboxy-1-methylethoxy)imino]acetyl]amino]-8-oxo-, trihydrochloride, (6R,7R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Double bond geometry as shown.



●3 HCl

REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 6 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2004:20537 CAPLUS Full-text

DOCUMENT NUMBER: 140:87699

TITLE: Remedies for diseases caused by vascular contraction or dilation

INVENTOR(S): Nakade, Shinji; Suzuki, Hidehiro; Habashita, Hiromu
 PATENT ASSIGNEE(S): Ono Pharmaceutical Co., Ltd., Japan
 SOURCE: PCT Int. Appl., 216 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: Japanese
 FAMILY ACC. NUM. COUNT: 2
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004002531	A1	20040108	WO 2003-JP8039	20030625 <--
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
AU 2003248245	A1	20040119	AU 2003-248245	20030625 <--
EP 1522314	A1	20050413	EP 2003-761797	20030625 <--
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK				
US 20060148844	A1	20060706	US 2005-519113	20051101 <--
PRIORITY APPLN. INFO.:			JP 2002-185546	A 20020626 <--
			WO 2003-JP8039	W 20030625 <--

OTHER SOURCE(S): MARPAT 140:87699

AB Remedies and/or preventives for diseases caused by vascular contraction or dilation which comprise EDG-5 regulators. EDG-5 regulators specifically bind to EDG-5 and show antagonism or agonism. Thus, an EDG-5 antagonist is useful in treating and/or preventing diseases caused by vascular contraction such as cerebrovascular spasmodic disease following subarachnoid hemorrhage or cerebral infarction, cardiovascular spasmodic disease, hypertension, kidney diseases, cardiac infarction, angina, arrhythmia, portal hypertension in association with cirrhosis and varicosity in association with cirrhosis. On the other hand, an EDG-5 agonist is useful in treating and/or preventing diseases caused by vascular dilation such as chronic headache (for example, hemicrania, tension headache, headache of the mixed type, cluster headache), hemorrhoid and cardiac diseases.

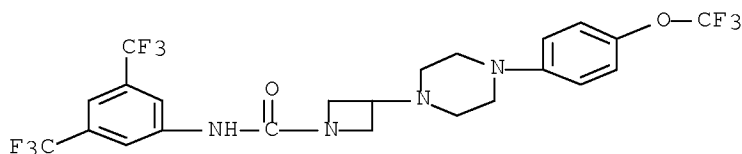
IT 642495-73-6P 642498-98-4P 642498-99-5P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

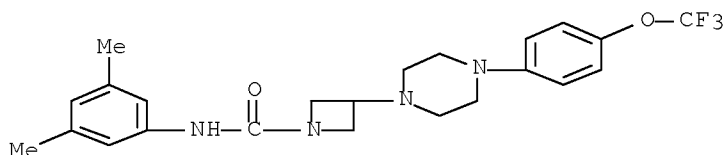
(EDG-5 agonists and antagonists as remedies for diseases caused by vascular contraction or dilation)

RN 642495-73-6 CAPLUS

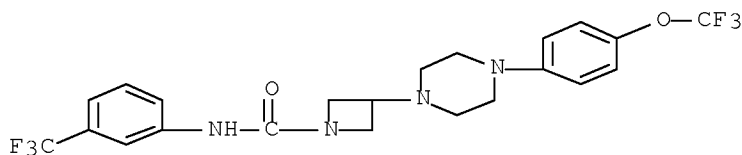
CN 1-Azetidinecarboxamide, N-[3,5-bis(trifluoromethyl)phenyl]-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



RN 642498-98-4 CAPLUS
 CN 1-Azetidinecarboxamide, N-(3,5-dimethylphenyl)-3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]- (CA INDEX NAME)



RN 642498-99-5 CAPLUS
 CN 1-Azetidinecarboxamide, 3-[4-[4-(trifluoromethoxy)phenyl]-1-piperazinyl]-N-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



REFERENCE COUNT: 15 THERE ARE 15 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 7 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2002:6377 CAPLUS Full-text

DOCUMENT NUMBER: 136:69695

TITLE: Preparation of β -lactam compounds as inhibitors of tryptase

INVENTOR(S): Bisacchi, Gregory S.; Slusarchyk, William A.; Treuner, Uwe; Sutton, James C.; Zahler, Robert; Seiler, Steven; Kronenthal, David R.; Randazzo, Michael E.; Schwinden, Mark D.; Xu, Zhongmin; Shi, Zhongping

PATENT ASSIGNEE(S): Bristol-Myers Squibb Co., USA

SOURCE: U.S., 171 pp., Cont.-in-part of U. S. Ser. No. 336,253, abandoned.

CODEN: USXXAM

DOCUMENT TYPE: Patent

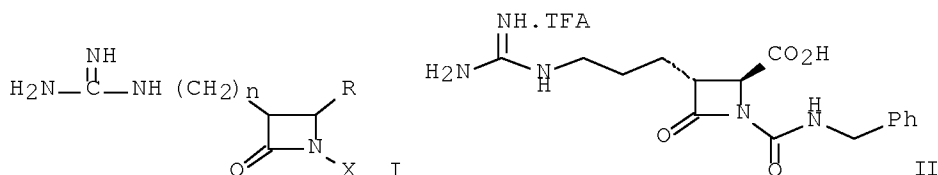
LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 6335324	B1	20020101	US 1999-458847	19991213 <--
PRIORITY APPLN. INFO.:			US 1998-90636P	P 19980625 <--
			US 1999-336253	B2 19990618 <--
OTHER SOURCE(S):	MARPAT	136:69695		

GI



AB Novel β -lactam compds., e.g. of formula I [R = CO₂H, alkoxy carbonyl, acyl, CO-heterocyclyl, etc.; X = acyl, CO-heterocyclyl, SO₂-alkyl, aminoalkylphenyl, etc.; n = 1-6], are prepared. These compds. inhibit tryptase as well as other enzyme systems or are selective tryptase inhibitors and are useful as antiinflammatory agents particularly in the treatment of chronic asthma (no data). Thus, II was prepared from (4S)-N-(tert-butyldimethylsilyl)azetidin-2-one-4-carboxylic acid, 1-chloro-3-iodopropane, N,N'-bis(benzyloxycarbonyl)-1-guanylpurazole and benzyl isocyanate.

IT 253173-23-8P 253173-24-9P

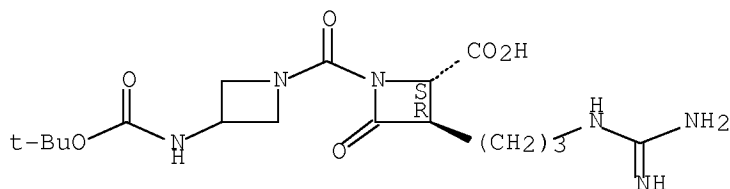
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of β -lactam compds. as inhibitors of tryptase)

RN 253173-23-8 CAPLUS

CN 2-Azetidinecarboxylic acid, 3-[3-[(aminoiminomethyl)amino]propyl]-1-[[3-[[[(1,1-dimethylethoxy)carbonyl]amino]-1-azetidinyl]carbonyl]-4-oxo-, monohydrochloride, (2S,3R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

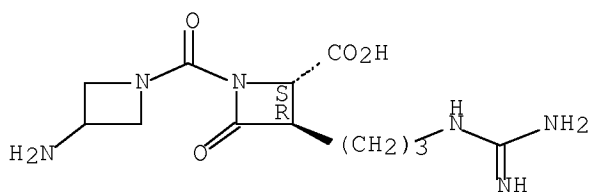


● HCl

RN 253173-24-9 CAPLUS

CN 2-Azetidinecarboxylic acid, 1-[(3-amino-1-azetidinyl)carbonyl]-3-[3-[(aminoiminomethyl)amino]propyl]-4-oxo-, dihydrochloride, (2S,3R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



● 2 HCl

IT 253176-37-3P 253176-39-5P

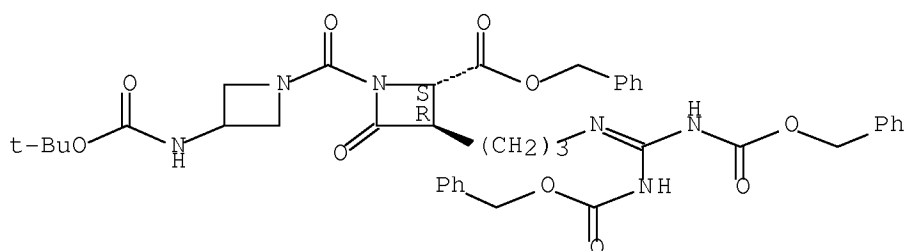
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of β -lactam compds. as inhibitors of tryptase)

RN 253176-37-3 CAPLUS

CN 2-Azetidinecarboxylic acid, 3-[3-[[bis[[(phenylmethoxy)carbonyl]amino]methylene]amino]propyl]-1-[[3-[[(1,1-dimethylethoxy)carbonyl]amino]-1-azetidiny]carbonyl]-4-oxo-, phenylmethyl ester, (2S,3R)- (CA INDEX NAME)

Absolute stereochemistry.



RN 253176-39-5 CAPLUS

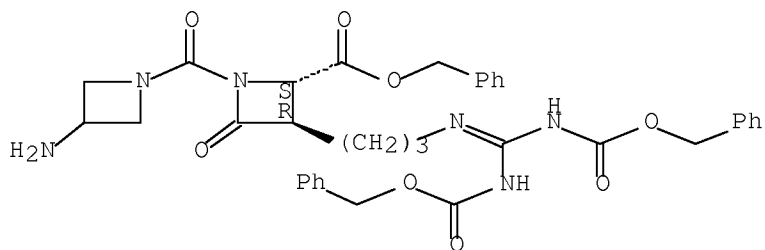
CN 2-Azetidinecarboxylic acid, 1-[(3-amino-1-azetidiny]carbonyl]-3-[3-[[bis[[(phenylmethoxy)carbonyl]amino]methylene]amino]propyl]-4-oxo-, phenylmethyl ester, (2S,3R)-, mono(trifluoroacetate) (9CI) (CA INDEX NAME)

CM 1

CRN 253176-38-4

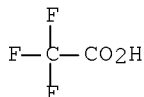
CMF C35 H38 N6 O8

Absolute stereochemistry.



CM 2

CRN 76-05-1
CMF C2 H F3 O2



REFERENCE COUNT: 26 THERE ARE 26 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 8 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2001:799296 CAPLUS Full-text

DOCUMENT NUMBER: 136:118095

TITLE: Structures of dinitroazetidine and three of its carbonyl derivatives

AUTHOR(S): Peiris, Suhithi M.; Gilardi, Richard; Dave, Paritosh R.; Russell, T. P.

CORPORATE SOURCE: Naval Research Lab, Washington, DC, 20375, USA

SOURCE: Journal of Chemical Crystallography (2001),

Volume Date 2000, 30(10), 647-653

CODEN: JCCYEV; ISSN: 1074-1542

PUBLISHER: Kluwer Academic/Plenum Publishers

DOCUMENT TYPE: Journal

LANGUAGE: English

AB The structures of 3,3-dinitroazetidine (DNAZ), DNAZ-COH, DNAZ-CO-DNAZ, and (DNAZ-CO)₂ have been determined DNAZ crystallizes in the space group P21/c with cell parameters a = 5.928(7) Å, b = 15.8380(2) Å, c = 12.6837(15) and β = 93.23°. DNAZ-COH crystallizes in the space group P21/c with cell parameters a = 10.3270(10) Å, b = 6.0615(12) Å, c = 11.2514(12), and β = 100.17°. DNAZ-CO-DNAZ crystallizes in the space group P21/n with cell parameters a = 10.3630(6) Å, b = 6.0901(5) Å, c = 19.9421(11) and β = 102.74°. (DNAZ-CO)₂ crystallizes in the orthorhombic space group Fdd2 with cell parameters a = 13.1956(10) Å, b = 19.524(3) Å, c = 10.3259(11). This last compound is unique in that the trans carbonyls are planar and the adjoining azetidine rings are also planar to within 0.1° in all torsion angles. DNAZ crystals have a dimeric asym. unit where two nonequal mols. are bonded to each other with a single hydrogen bond. Like other four-membered rings with gem-dinitro groups, most of these dinitroazetidines display mutually perpendicular nitro groups.

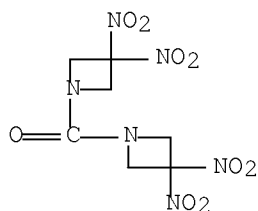
IT 211429-18-4, Azetidine, 1,1'-carbonylbis[3,3-dinitro-

RL: PRP (Properties)

(crystal structure; crystal and mol. structure of dinitroazetidine and carbonyl derivs.)

RN 211429-18-4 CAPLUS

CN Azetidine, 1,1'-carbonylbis[3,3-dinitro- (9CI) (CA INDEX NAME)

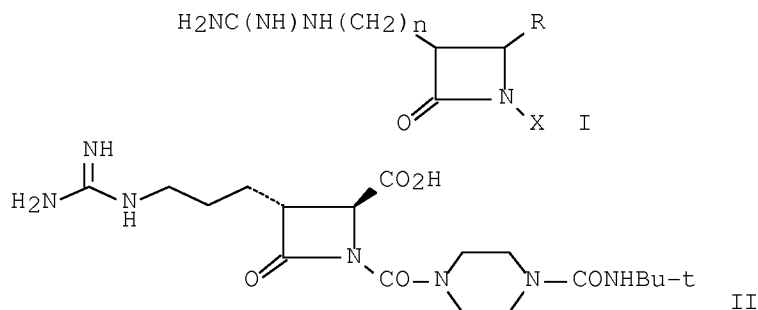


REFERENCE COUNT: 19 THERE ARE 19 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 9 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN
 ACCESSION NUMBER: 1999:819347 CAPLUS Full-text
 DOCUMENT NUMBER: 132:64103
 TITLE: Preparation of amidino and guanidino azetidinone compounds as tryptase inhibitors
 INVENTOR(S): Bisacchi, Gregory; Slusarchyk, William A.; Treuner, Uwe; Sutton, James C.; Zahler, Robert; Seiler, Steven; Kronenthal, David R.; Randazzo, Michael E.; Xu, Zhongmin; Shi, Zhongping; Schwinden, Mark D.
 PATENT ASSIGNEE(S): Bristol-Myers Squibb Co., USA
 SOURCE: PCT Int. Appl., 326 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 2
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9967215	A1	19991229	WO 1999-US13811	19990618 <--
W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, UZ, VN, YU, ZA, ZW				
RW: GH, GM, KE, LS, MW, SD, SL, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
CA 2336003	A1	19991229	CA 1999-2336003	19990618 <--
AU 9946950	A	20000110	AU 1999-46950	19990618 <--
AU 752320	B2	20020912		
EP 1089973	A1	20010411	EP 1999-930402	19990618 <--
EP 1089973	B1	20051109		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, FI				
TR 200003859	T2	20010723	TR 2000-3859	19990618 <--
BR 9911373	A	20010918	BR 1999-11373	19990618 <--
HU 2001004308	A2	20020429	HU 2001-4308	19990618 <--
HU 2001004308	A3	20020628		
JP 2002518478	T	20020625	JP 2000-555869	19990618 <--
RU 2211832	C2	20030910	RU 2001-102266	19990618 <--
NZ 507627	A	20031219	NZ 1999-507627	19990618 <--
AT 309211	T	20051115	AT 1999-930402	19990618 <--
ES 2249017	T3	20060316	ES 1999-930402	19990618 <--
TW 548270	B	20030821	TW 1999-88110361	19990621 <--
ZA 2000006028	A	20020725	ZA 2000-6028	20001025 <--

MX 2000PA11532	A	20010629	MX 2000-PA11532	20001123 <--
NO 2000006380	A	20001214	NO 2000-6380	20001214 <--
IN 2003MN00909	A	20050429	IN 2003-MN909	20030924 <--
PRIORITY APPLN. INFO.:			US 1998-90636P	P 19980625 <--
			WO 1999-US13811	W 19990618 <--
OTHER SOURCE(S):		MARPAT 132:64103		
GI				



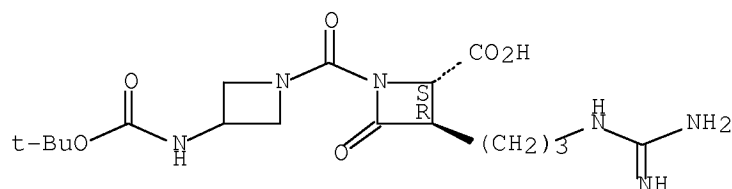
AB Novel β -lactam compds., e.g. of formula I [R - CO₂H, CONH-alkyl, etc.; X = CONH(CH₂)₂NHCO₂alkyl, etc.; n = 1-6;], are prepared as inhibitors of in vivo enzyme systems including tryptase, thrombin, trypsin, factor Xa, factor VIIa, and urokinase-type plasminogen activator (no data). The tryptase activity makes the title compds. useful as antiinflammatory agents in the treatment of chronic asthma and allergic rhinitis. Thus, II was prepared from (4S)-N-(tert-butyltrimethylsilyl)azetidin-2-one-4-carboxylic acid, tert-butyl-1-piperazine carboxylate and tert-Bu isocyanate.

IT 253173-23-8P 253173-24-9P
 RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
 (preparation of amidino and guanidino azetidinone compds. as tryptase inhibitors)

RN 253173-23-8 CAPLUS

CN 2-Azetidinecarboxylic acid, 3-[3-[(aminoiminomethyl)amino]propyl]-1-[[3-[[[1,1-dimethylethoxy]carbonyl]amino]-1-azetidiny]carbonyl]-4-oxo-, monohydrochloride, (2S,3R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



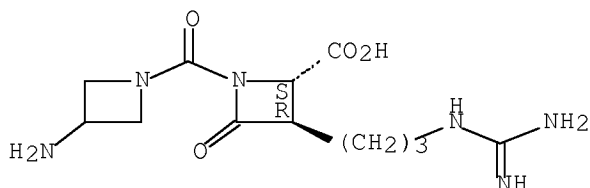
● HCl

RN 253173-24-9 CAPLUS

CN 2-Azetidinecarboxylic acid, 1-[(3-amino-1-azetidiny]carbonyl]-3-[3-

[(aminoiminomethyl)amino]propyl]-4-oxo-, dihydrochloride, (2S,3R)- (9CI)
(CA INDEX NAME)

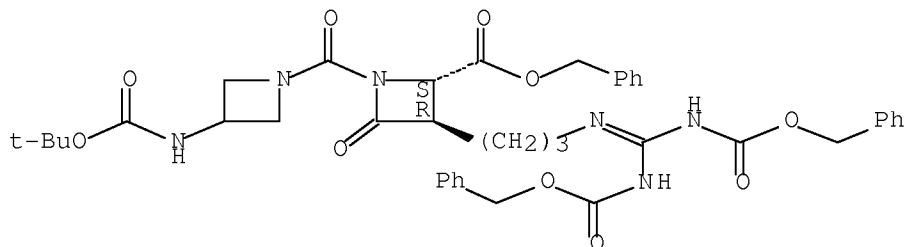
Absolute stereochemistry.



● 2 HCl

IT 253176-37-3P 253176-39-5P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
(Reactant or reagent)
(preparation of amidino and guanidino azetidinone compds. as tryptase
inhibitors)
RN 253176-37-3 CAPLUS
CN 2-Azetidinecarboxylic acid, 3-[3-[[bis[[(phenylmethoxy)carbonyl]amino]meth-
ylene]amino]propyl]-1-[[3-[[(1,1-dimethylethoxy)carbonyl]amino]-1-
azetidiny]l]carbonyl]-4-oxo-, phenylmethyl ester, (2S,3R)- (CA INDEX NAME)

Absolute stereochemistry.

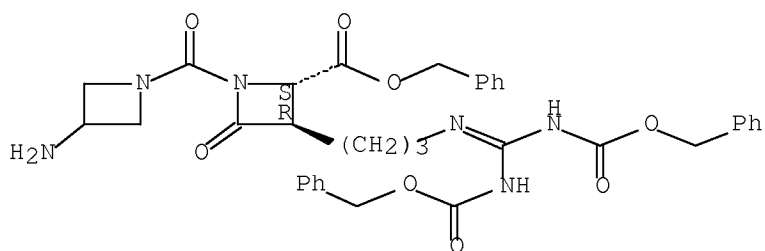


RN 253176-39-5 CAPLUS
CN 2-Azetidinecarboxylic acid, 1-[(3-amino-1-azetidiny]l)carbonyl]-3-[3-
[[bis[[(phenylmethoxy)carbonyl]amino]methylene]amino]propyl]-4-oxo-,
phenylmethyl ester, (2S,3R)-, mono(trifluoroacetate) (9CI) (CA INDEX
NAME)

CM 1

CRN 253176-38-4
CMF C35 H38 N6 O8

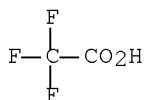
Absolute stereochemistry.



CM 2

CRN 76-05-1

CMF C2 H F3 O2



REFERENCE COUNT: 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 10 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 1999:411860 CAPLUS Full-text

DOCUMENT NUMBER: 131:144472

TITLE: Preparation of 1-substituted 3,3-dinitroazetidines

AUTHOR(S): Hiskey, Michael A.; Johnson, M. Catherine; Chavez, David E.

CORPORATE SOURCE: Los Alamos National Laboratory, Los Alamos, NM, 87545, USA

SOURCE: Journal of Energetic Materials (1999), 17(2 & 3), 233-252

CODEN: JOEMDK; ISSN: 0737-0652

PUBLISHER: Dowden, Brodman & Devine, Inc.

DOCUMENT TYPE: Journal

LANGUAGE: English

AB DNAZ (3,3-dinitroazetidine) readily undergoes nucleophilic addition reactions with electron deficient haloaroms., cyanogen bromide, S-methyl-N-nitroisothiourea and dichloroglyoxime to afford novel, energetic 1-substituted 3,3-dinitroazetidines. N-Nitrosation of DNAZ readily occurs, and Mannich condensations of DNAZ with polynitro alcs. are also described. The synthesis and properties of the energetic materials derived from these reactions are discussed.

IT 236102-62-8P 236102-63-9P 236102-65-1P
236102-66-2P

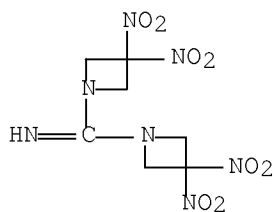
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(preparation and explosive properties of 1-substituted 3,3-dinitroazetidines)

RN 236102-62-8 CAPLUS

CN Azetidine, 1,1'-carbonimidoylbis-, mononitrate (9CI) (CA INDEX NAME)

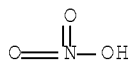
CM 1

CRN 236102-61-7
CMF C7 H9 N7 O8



CM 2

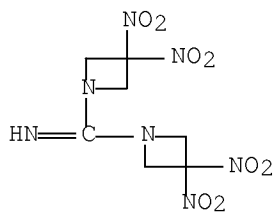
CRN 7697-37-2
CMF H N O3



RN 236102-63-9 CAPLUS
CN Azetidine, 1,1'-carbonimidoylbis-, compd. with nitronitramide (1:1) (9CI)
(CA INDEX NAME)

CM 1

CRN 236102-61-7
CMF C7 H9 N7 O8



CM 2

CRN 114045-20-4
CMF H N3 O4

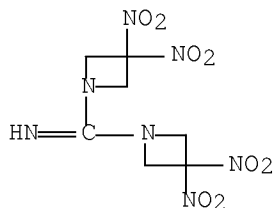


RN 236102-65-1 CAPLUS
CN Azetidine, 1,1'-carbonimidoylbis-, monoperoxylate (9CI) (CA INDEX NAME)

CM 1

CRN 236102-61-7

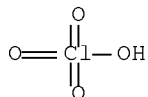
CMF C7 H9 N7 O8



CM 2

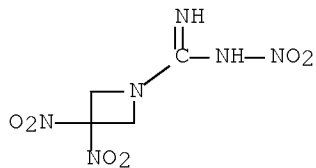
CRN 7601-90-3

CMF Cl H O4



RN 236102-66-2 CAPLUS

CN 1-Azetidinecarboximidamide, N,3,3-trinitro- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 20 THERE ARE 20 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 11 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 1999:3288 CAPLUS Full-text

DOCUMENT NUMBER: 130:66390

TITLE: Preparation of 1-benzenesulfonyl-1,3-dihydroindol-2-ones as vasopressin and/or oxytocin antagonists

INVENTOR(S): Di Malta, Alain; Foulon, Loic; Garcia, Georges; Nisato, Dino; Roux, Richard; Serradeil-Legal, Claudine; Valette, Gerard; Wagnon, Jean

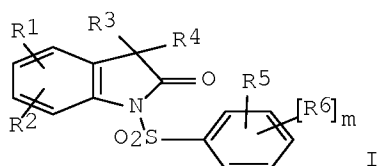
PATENT ASSIGNEE(S): Sanofi, Fr.

SOURCE: U.S., 53 pp., Cont.-in-part of U.S. Ser. No. 129,310, abandoned.

CODEN: USXXAM

DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 3
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 5849780	A	19981215	US 1994-323921	19941017 <--
FR 2686878	A1	19930806	FR 1992-1034	19920130 <--
FR 2686878	B1	19950630		
FR 2708605	A1	19950210	FR 1993-9404	19930730 <--
EP 636608	A1	19950201	EP 1994-401737	19940728 <--
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LI, LU, MC, NL, PT, SE				
US 5663431	A	19970902	US 1995-477571	19950607 <--
US 5686624	A	19971111	US 1995-473302	19950607 <--
US 5728723	A	19980317	US 1995-478738	19950607 <--
US 5726322	A	19980310	US 1997-824305	19970326 <--
PRIORITY APPLN. INFO.:			FR 1992-1034	A 19920130 <--
			FR 1993-9404	A 19930730 <--
			US 1993-129310	B2 19930930 <--
			EP 1994-401737	A 19940728 <--
			US 1994-323921	A3 19941017 <--
			US 1995-473302	A3 19950607 <--
OTHER SOURCE(S):			MARPAT 130:66390	
GI				

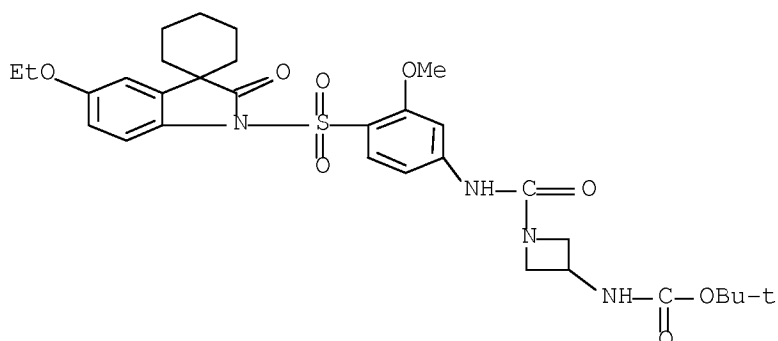


AB The title compds. [I; R1, R2 = H, OH, halo, etc.; R3R4 together with the carbon to which they are bonded = an optionally fused, (un)saturated (un)substituted C3-12 hydrocarbon ring; R5, R6 = H, halo, C1-7 alkyl, etc.; m = 1-4], having an affinity for the vasopressin V1 and V2 and/or oxytocin receptors, were prepared Thus, treatment of 5-chloro-1,3-dihydro-3-spirocyclohexaneindol-2-one with NaH in THF followed by addition of 2-methoxy-4-nitrobenzenesulfonyl chloride afforded I [R1 = 5-Cl; R2 = H; R3R4 = (CH2)5; R5 = 2-MeO; R6 = 4-NO2]. Biol. data for compds. I are given.

IT 217806-19-4P
 RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
 (preparation of 1-benzenesulfonyl-1,3-dihydroindol-2-ones as vasopressin and/or oxytocin antagonists)

RN 217806-19-4 CAPLUS

CN Carbamic acid, [1-[[[4-[(5'-ethoxy-2'-oxospiro[cyclohexane-1,3'-[3H]indol]-1'(2'H)-yl)sulfonyl]-3-methoxyphenyl]amino]carbonyl]-3-azetidiny]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



REFERENCE COUNT: 17 THERE ARE 17 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L17 ANSWER 12 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 1998:460439 CAPLUS Full-text

DOCUMENT NUMBER: 129:177578

TITLE: Condensation reaction of 3,3-dinitroazetidine

AUTHOR(S): Zhang, Jiaoqiang; Zhu, Chunhua; Zhang, Gao

CORPORATE SOURCE: Xi'an Modern Chem. Res. Inst., Xi'an, 710065, Peop. Rep. China

SOURCE: Hanneng Cailiao (1998), 6(2), 73-77

CODEN: HACAFQ; ISSN: 1006-9941

PUBLISHER: Hanneng Cailiao Bianjibu

DOCUMENT TYPE: Journal

LANGUAGE: Chinese

AB The condensation of 3,3-dinitroazetidine with phosgene and oxalyl chloride resulted in the isolation of two tetranitro diamides. The thermal decomposition kinetics of the two tetranitro diamides were determined by differential scanning calorimetry, and their structures were confirmed by IR and NMR spectroscopy. Quantum chemical calcns. for bis(3,3-dinitroazetidine)one (from reaction with phosgene) were carried out using the SCF-MO-AM1 method, and an optimal geometric configuration and electronic structure were obtained.

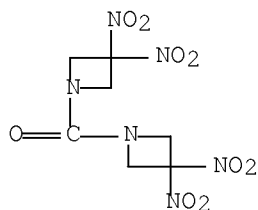
IT 211429-18-4F

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)

(synthesis and properties of tetranitro diamides prepared by amidation of 3,3-dinitroazetidine with phosgene or oxalyl chloride)

RN 211429-18-4 CAPLUS

CN Azetidine, 1,1'-carbonylbis[3,3-dinitro- (9CI) (CA INDEX NAME)



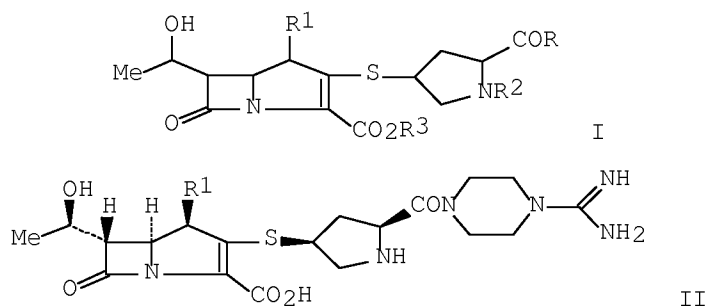
L17 ANSWER 13 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 1994:269926 CAPLUS Full-text

DOCUMENT NUMBER: 120:269926

TITLE: Preparation of 2-(2-carbamoyl-4-pyrrolidinylthio)carbapenems as antibiotics
 INVENTOR(S): Kawamoto, Isao; Endo, Rokuro; Miyauchi, Masao; Ishikawa, Katsuya; Watanabe, Katsuhiko; Yasuda, Hiroshi; Ohya, Satoshi; Utsui, Yukio
 PATENT ASSIGNEE(S): Sankyo Co., Ltd., Japan
 SOURCE: Eur. Pat. Appl., 139 pp.
 CODEN: EPXXDW
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 2
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 560613	A1	19930915	EP 1993-301872	19930311 <--
EP 560613	B1	19981125		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LI, LU, MC, NL, PT, SE				
HU 64071	A2	19931129	HU 1993-677	19930310 <--
HU 218676	B	20001028		
JP 06041131	A	19940215	JP 1993-49082	19930310 <--
FI 103046	B	19990415	FI 1993-1079	19930310 <--
FI 103046	B1	19990415		
CZ 290002	B6	20020515	CZ 1993-393	19930310 <--
CA 2091486	A1	19930912	CA 1993-2091486	19930311 <--
NO 9300884	A	19930913	NO 1993-884	19930311 <--
NO 303730	B1	19980824		
AU 9335151	A	19930916	AU 1993-35151	19930311 <--
AU 664422	B2	19951116		
ZA 9301761	A	19930928	ZA 1993-1761	19930311 <--
CN 1079472	A	19931215	CN 1993-104412	19930311 <--
CN 1037180	B	19980128		
IL 105020	A	19970415	IL 1993-105020	19930311 <--
RU 2097383	C1	19971127	RU 1993-4678	19930311 <--
AT 173733	T	19981215	AT 1993-301872	19930311 <--
ES 2125945	T3	19990316	ES 1993-301872	19930311 <--
JP 06172356	A	19940621	JP 1993-227141	19930913 <--
JP 3344662	B2	20021111		
JP 06199860	A	19940719	JP 1993-228457	19930914 <--
US 5866564	A	19990202	US 1997-889542	19970708 <--
PRIORITY APPLN. INFO.:				
			JP 1992-52163	A 19920311 <--
			JP 1992-244953	A 19920914 <--
			JP 1992-246578	A 19920916 <--
			JP 1991-131545	A 19910604 <--
			JP 1991-345737	A 19911227 <--
			JP 1992-30521	A 19920218 <--
			JP 1992-91283	A 19920410 <--
			US 1992-894004	B1 19920603 <--
			US 1993-29779	B1 19930311 <--
			US 1993-81848	B1 19930622 <--
			US 1994-288987	B2 19940811 <--
			US 1994-293378	B2 19940819 <--
			US 1995-472850	A3 19950606 <--
OTHER SOURCE(S): MARPAT 120:269926				
GI				



AB Title compds. [I; R = substituted amino, N-attached heterocyclyl; R1 = H, Me; R2 = H, alk(en)yl, alkynyl, etc.; R3 = H, carboxy-protective group, etc.] were prepared Thus, 4-nitrobenzyl (1R,5S,6S)-6-[(1R)-1-hydroxyethyl]- 1-methyl-2-oxo-1-carbapenam-3-carboxylate was condensed with (2S, 4S)-4-mercapto-2-[4-(4-nitrobenzyloxycarbonylamidino)piperazin-1-ylcarbonyl]-1-(4-nitrobenzyloxycarbonyl)pyrrolidine (preparation given) to give title compound II which had MIC of ≤ 0.01 (units not given) against Staphylococcus aureus 209P and Escherichia coli NIHJ.

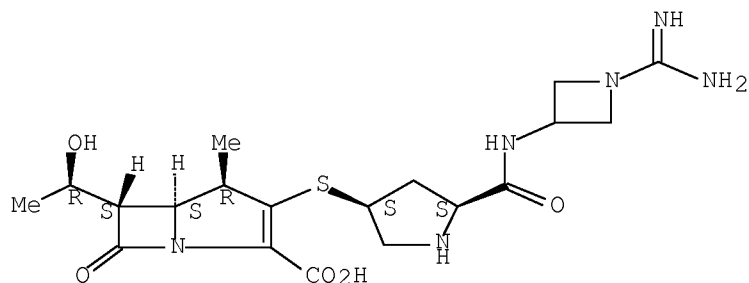
IT 153835-94-0P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)
(preparation of, as antibiotic)

RN 153835-94-0 CAPLUS

CN 1-Azabicyclo[3.2.0]hept-2-ene-2-carboxylic acid, 3-[[5-[[[1-(aminoiminomethyl)-3-azetidiny]amino]carbonyl]-3-pyrrolidinyl]thio]-6-(1-hydroxyethyl)-4-methyl-7-oxo-, [4R-[3(3S*,5S*),4 α ,5 β ,6 β (R*)]]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



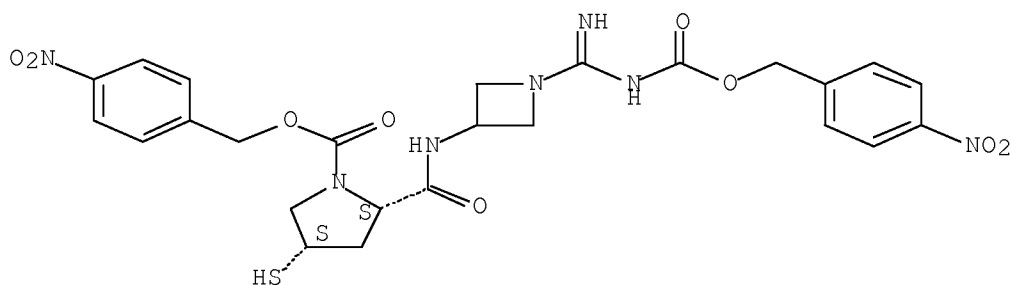
IT 153836-82-9

RL: RCT (Reactant); RACT (Reactant or reagent)
(reaction of, in preparation of antibiotic)

RN 153836-82-9 CAPLUS

CN 1-Pyrrolidinecarboxylic acid, 2-[[[1-[imino[[[(4-nitrophenyl)methoxy]carbonyl]amino]methyl]-3-azetidiny]amino]carbonyl]-4-mercapto-, (4-nitrophenyl)methyl ester, (2S-cis)- (9CI) (CA INDEX NAME)

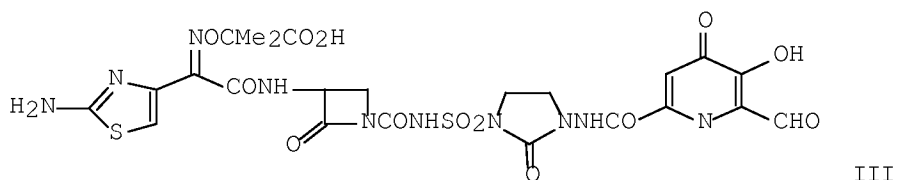
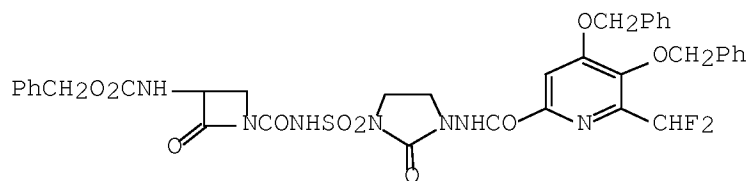
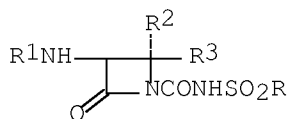
Absolute stereochemistry.



L17 ANSWER 14 OF 15 CAPLUS COPYRIGHT 2008 ACS on STN
 ACCESSION NUMBER: 1990:440323 CAPLUS Full-text
 DOCUMENT NUMBER: 113:40323
 TITLE: Preparation of 3-acylamino-1-
 [[(heterocyclylsulfonyl)amino]carbonyl]-2-azetidinones
 as antibiotics
 INVENTOR(S): Treuner, Uwe D.
 PATENT ASSIGNEE(S): E. R. Squibb and Sons, Inc., USA
 SOURCE: Eur. Pat. Appl., 69 pp.
 CODEN: EPXXDW
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 336369	A1	19891011	EP 1989-105883	19890404 <--
R: AT, BE, CH, DE, ES, FR, GB, GR, IT, LI, LU, NL, SE				
JP 01305081	A	19891208	JP 1989-86701	19890404 <--
PRIORITY APPLN. INFO.:			US 1988-177207	A 19880404 <--
OTHER SOURCE(S):	MARPAT 113:40323			

GI



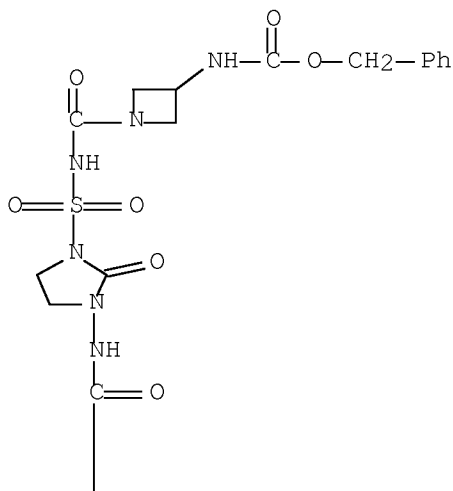
AB The title compds. [I; R = [(oxopyridinylcarbonyl)amino]oxoimidazolo, -
dioxopiperazino, etc.; R1 = acyl; R2, R3 = H, alkyl, alkenyl, etc.] were
prepared as antibiotics (no data). Thus, I.CF3CO2H [R = [(tert-
butoxycarbonyl)amino]oxoimidazolo] (preparation given) was treated with
CF3CONMeSiMe3 and the product condensed with 1H-benzotriazol-1-yl 6-
difluoromethyl-4,5-bis(phenylmethoxy)-2-pyridinecarboxylate (preparation
given) to give, after neutralization, title compound II.Na which was converted
in 2 steps to title compound (3S)-(Z)-III.2Na.

IT 127235-37-4P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
(Reactant or reagent)
(preparation and reaction of, in preparation of antibiotics)

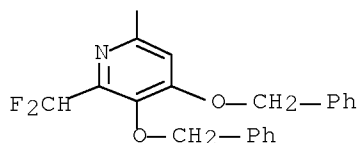
RN 127235-37-4 CAPLUS

CN Carbamic acid, [1-[[[3-[[[6-(difluoromethyl)-4,5-bis(phenylmethoxy)-2-
pyridinyl]carbonyl]amino]-2-oxo-1-imidazolidinyl]sulfonyl]amino]carbonyl]-
3-azetidiny]-, phenylmethyl ester (9CI) (CA INDEX NAME)

PAGE 1-A



PAGE 2-A



TITLE: Preparation of N-substituted azetidinone derivatives
as antibiotics
PATENT ASSIGNEE(S): E. R. Squibb and Sons, Inc., USA
SOURCE: Belg., 173 pp.
CODEN: BEXXAL
DOCUMENT TYPE: Patent
LANGUAGE: French
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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BE 905502	A1	19870116	BE 1986-217217	19860926 <--
US 4743685	A	19880510	US 1986-907441	19860915 <--
DK 8604603	A	19870327	DK 1986-4603	19860926 <--
FI 8603890	A	19870327	FI 1986-3890	19860926 <--
FR 2587700	A1	19870327	FR 1986-13466	19860926 <--
FR 2587700	B1	19880916		
NO 8603837	A	19870327	NO 1986-3837	19860926 <--
SE 8604089	A	19870327	SE 1986-4089	19860926 <--
AU 8663156	A	19870402	AU 1986-63156	19860926 <--
AU 600912	B2	19900830		
DE 3632876	A1	19870402	DE 1986-3632876	19860926 <--
CN 86106980	A	19870408	CN 1986-106980	19860926 <--
GB 2181130	A	19870415	GB 1986-23151	19860926 <--
GB 2181130	B	19890920		
NL 8602446	A	19870416	NL 1986-2446	19860926 <--
JP 62084082	A	19870417	JP 1986-229474	19860926 <--
ZA 8607373	A	19870527	ZA 1986-7373	19860926 <--
DD 250121	A5	19870930	DD 1986-295013	19860926 <--
HU 44535	A2	19880328	HU 1986-4124	19860926 <--
HU 198044	B	19890728		
ES 2001995	A6	19880701	ES 1986-2204	19860926 <--
CH 670828	A5	19890714	CH 1986-3860	19860926 <--
PL 151546	B1	19900928	PL 1986-261575	19860926 <--
IL 80160	A	19910415	IL 1986-80160	19860926 <--
CA 1308719	C	19921013	CA 1986-519170	19860926 <--
PRIORITY APPLN. INFO.:			US 1985-780479	A 19850926 <--
OTHER SOURCE(S):		CASREACT 108:55763; MARPAT 108:55763		
GI				

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The title compds. I [R = Q1, Q2, etc., wherein A1 = single bond, NHCO, NH, NHHNCO, A2 = single bond, NH, (CH₂)₂NH, CONHNH, A3 = single bond, CH:CH, (CH₂)_n (n = 1-4); R1 = acyl; R2, R3 = H, alkyl, alkenyl, alkynyl, cycloalkyl, (un)substituted Ph, heterocyclyl, or one of R2 and R3 is H, the other is N3, halomethyl, dihalomethyl, alkoxy carbonyl, 2-phenylethenyl, etc.], useful as antibiotics (no data), were prepared Azetidinone derivative (S)-II (R1 = PhCH₂OCO, R2 = R3 = H) was treated with ClSO₂NCO to give (chlorosulfonyl)aminocarbonylazetidinone derivative (S)-I (R1 = PhCH₂OCO, R2 = R3 = H, R = Cl) (III). To a solution of III in EtOAc was added a mixture of 1,4-dihydro-5-hydroxy-4-oxo-N-(2-oxo-1-imidazolidinyl)-2- pyridinecarboxamide and CF₃CON(SiMe₃)Me in EtOAc. Following addition of CH₂Cl₂, the mixture was stirred overnight at room temperature to give the corresponding imidazolidinylsulfonylaminocarbonylazetidinone derivative, which was deprotected with PhSMc/CF₃CO₂H and subsequently acylated with (Z)-2-amino- α -

[[2-(diphenylmethoxy)-1,1-dimethyl-2-oxoethoxy]imino]- 4-thiazoleacetic acid to give IV (R4 = Ph2CH). IV was then deprotected with PhOMe/CF3CO2H to give, after workup, azetidinone derivative IV.2Na (R4 = H).

IT 112334-33-5P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation and reaction of, in synthesis of antibiotic)

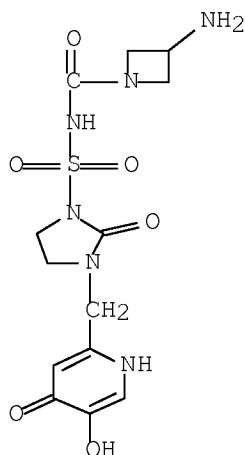
RN 112334-33-5 CAPLUS

CN 1-Azetidinecarboxamide, 3-amino-N-[[3-[(1,4-dihydro-5-hydroxy-4-oxo-2-pyridinyl)methyl]-2-oxo-1-imidazolidinyl]sulfonyl]-, mono(trifluoroacetate) (salt) (9CI) (CA INDEX NAME)

CM 1

CRN 112334-32-4

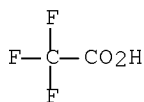
CMF C13 H18 N6 O6 S



CM 2

CRN 76-05-1

CMF C2 H F3 O2



=> fil stng

COST IN U.S. DOLLARS

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
97.83	331.91

SINCE FILE	TOTAL
ENTRY	SESSION

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	ENTRY	SESSION
FULL ESTIMATED COST	0.84	332.75
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-12.00

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